



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 09:16 PM EDT

PDB ID : 5X19 / pdb_00005x19
Title : CO bound cytochrome c oxidase at 100 micro sec after pump laser irradiation to release CO from O2 reduction center
Authors : Shimada, A.; Kubo, M.; Baba, S.; Yamashita, K.; Hirata, K.; Ueno, G.; Nomura, T.; Kimura, T.; Shinzawa-Itoh, K.; Baba, J.; Hatano, K.; Eto, Y.; Miyamoto, A.; Murakami, H.; Kumasaka, T.; Owada, S.; Tono, K.; Yabashi, M.; Yamaguchi, Y.; Yanagisawa, S.; Sakaguchi, M.; Ogura, T.; Komiya, R.; Yan, J.; Yamashita, E.; Yamamoto, M.; Ago, H.; Yoshikawa, S.; Tsukihara, T.
Deposited on : 2017-01-25
Resolution : 2.20 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	: 4-5-2 with Phenix2.0
Mogul	: 2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	: 2.0
EDS	: 3.0
Buster-report	: wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	: 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	: 9.0.010 (Gargrove)
Density-Fitness	: 1.0.12

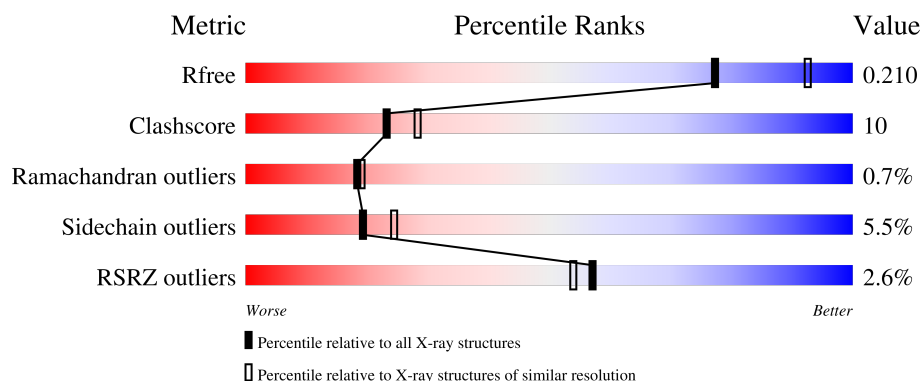
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

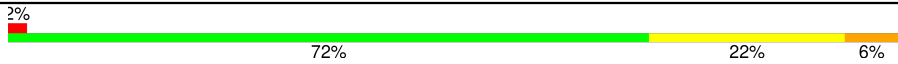
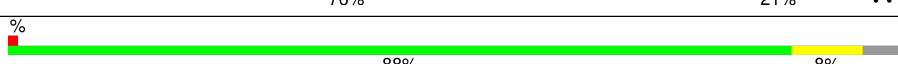
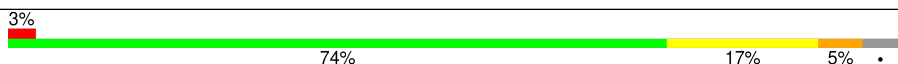
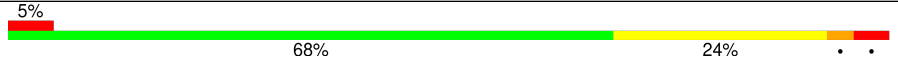
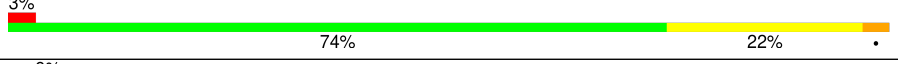
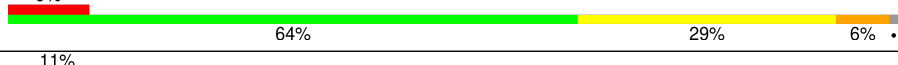
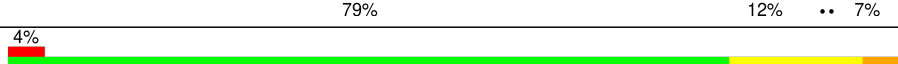
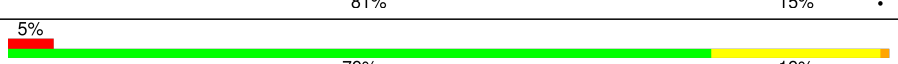
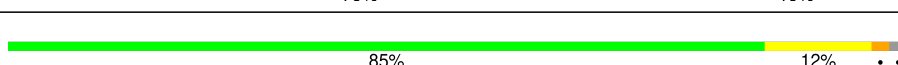
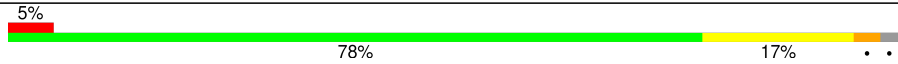
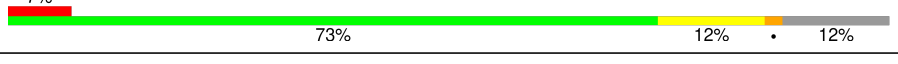
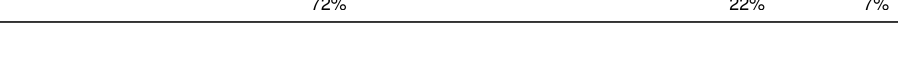
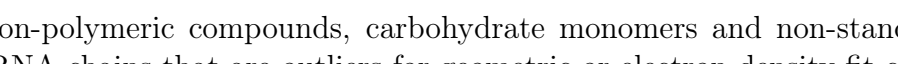
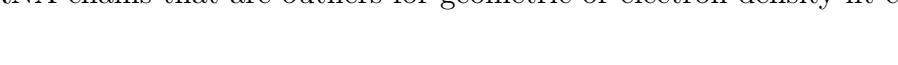
The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	514	
1	N	514	
2	B	227	

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Ideal geometry (proteins) : Engh & Huber (2001)
 Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
 Validation Pipeline (wwPDB-VP) : 2.49

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Mol	Chain	Length	Quality of chain
2	O	227	
3	C	261	
3	P	261	
4	D	147	
4	Q	147	
5	E	109	
5	R	109	
6	F	98	
6	S	98	
7	G	85	
7	T	85	
8	H	85	
8	U	85	
9	I	73	
9	V	73	
10	J	59	
10	W	59	
11	K	56	
11	X	56	
12	L	47	
12	Y	47	
13	M	46	
13	Z	46	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
18	CMO	N	606	-	-	X	-
21	EDO	A	612	-	-	X	-
21	EDO	A	613	-	-	X	-
25	UNX	C	301	-	-	-	X
25	UNX	P	302	-	-	-	X
26	CDL	T	104	-	-	X	-

2 Entry composition

There are 30 unique types of molecules in this entry. The entry contains 32242 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Cytochrome c oxidase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	514	Total	C	N	O	S	0	7	0
			4068	2718	628	686	36			
1	N	514	Total	C	N	O	S	0	0	0
			4027	2691	623	678	35			

- Molecule 2 is a protein called Cytochrome c oxidase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	227	Total	C	N	O	S	0	0	0
			1824	1185	281	340	18			
2	O	227	Total	C	N	O	S	0	0	0
			1824	1185	281	340	18			

- Molecule 3 is a protein called Cytochrome c oxidase subunit 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	259	Total	C	N	O	S	0	0	0
			2110	1412	336	350	12			
3	P	259	Total	C	N	O	S	0	0	0
			2110	1412	336	350	12			

- Molecule 4 is a protein called Cytochrome c oxidase subunit 4 isoform 1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	144	Total	C	N	O	S	0	0	0
			1195	777	196	218	4			
4	Q	144	Total	C	N	O	S	0	0	0
			1195	777	196	218	4			

- Molecule 5 is a protein called Cytochrome c oxidase subunit 5A, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	E	105	Total	C	N	O	S	0	0	0
			852	544	144	162	2			
5	R	105	Total	C	N	O	S	0	0	0
			852	544	144	162	2			

- Molecule 6 is a protein called Cytochrome c oxidase subunit 5B, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	F	98	Total	C	N	O	S	0	0	0
			748	464	134	145	5			
6	S	98	Total	C	N	O	S	0	0	0
			748	464	134	145	5			

- Molecule 7 is a protein called Cytochrome c oxidase subunit 6A2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
7	G	84	Total 675	C 431	N 129	O 113	P 1	S 1	0	0	0
7	T	84	Total 675	C 431	N 129	O 113	P 1	S 1	0	0	0

- Molecule 8 is a protein called Cytochrome c oxidase subunit 6B1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
8	H	79	Total	C	N	O	S	0	0	0
			662	417	121	119	5			
8	U	79	Total	C	N	O	S	0	0	0
			662	417	121	119	5			

- Molecule 9 is a protein called Cytochrome c oxidase subunit 6C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
9	I	73	Total	C	N	O	S	0	1	0
			609	398	107	100	4			
9	V	73	Total	C	N	O	S	0	2	0
			617	406	107	100	4			

- Molecule 10 is a protein called Cytochrome c oxidase subunit 7A1, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	J	58	Total	C	N	O	S	0	0	0
			460	297	78	82	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
10	W	58	Total	C	N	O	S	0	0	0
			460	297	78	82	3			

- Molecule 11 is a protein called Cytochrome c oxidase subunit 7B, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
11	K	49	Total	C	N	O	S	0	0	0
			384	250	65	67	2			
11	X	49	Total	C	N	O	S	0	0	0
			384	250	65	67	2			

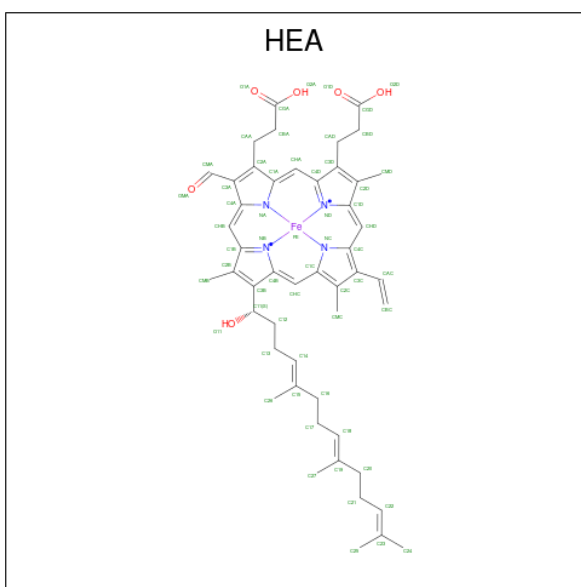
- Molecule 12 is a protein called Cytochrome c oxidase subunit 7C, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
12	L	46	Total	C	N	O	S	0	0	0
			380	254	64	60	2			
12	Y	46	Total	C	N	O	S	0	0	0
			380	254	64	60	2			

- Molecule 13 is a protein called Cytochrome c oxidase subunit 8B, mitochondrial.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	M	43	Total	C	N	O	0	0	0
			335	223	53	59			
13	Z	43	Total	C	N	O	0	0	0
			335	223	53	59			

- Molecule 14 is HEME-A (CCD ID: HEA) (formula: $C_{49}H_{56}FeN_4O_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
14	A	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	A	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	N	1	Total 60	C 49	Fe 1	N 4	O 6	0	0
14	N	1	Total 60	C 49	Fe 1	N 4	O 6	0	0

- Molecule 15 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
15	A	1	Total	Cu	0	0
			1	1		
15	N	1	Total	Cu	0	0
			1	1		

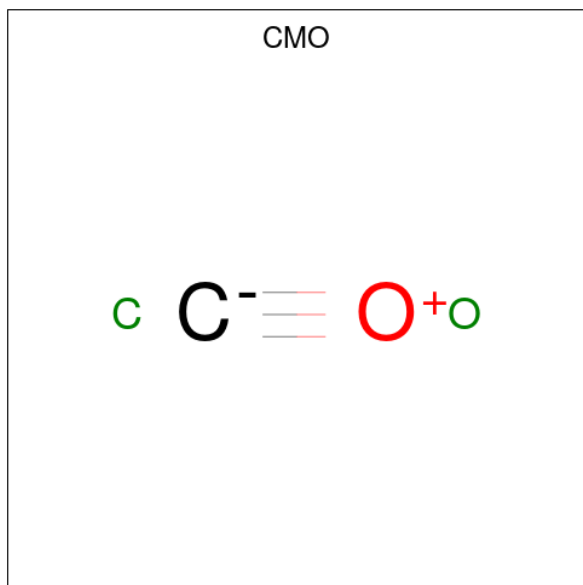
- Molecule 16 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
16	A	1	Total	Mg	0	0
			1	1		
16	N	1	Total	Mg	0	0
			1	1		

- Molecule 17 is SODIUM ION (CCD ID: NA) (formula: Na).

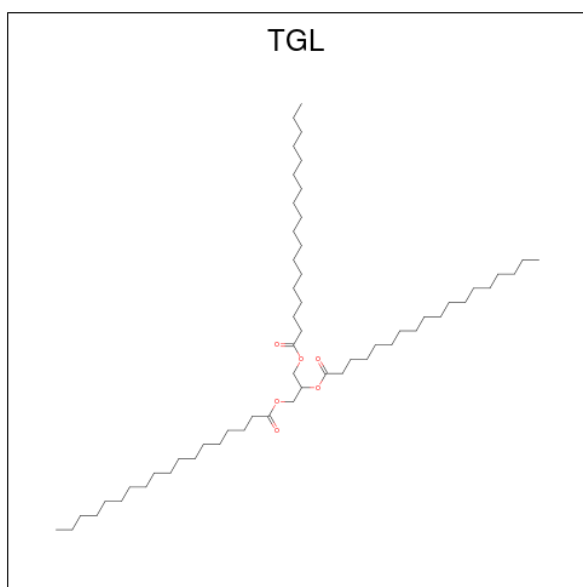
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
17	A	1	Total	Na	0	0
			1	1		
17	N	1	Total	Na	0	0
			1	1		

- Molecule 18 is CARBON MONOXIDE (CCD ID: CMO) (formula: CO).



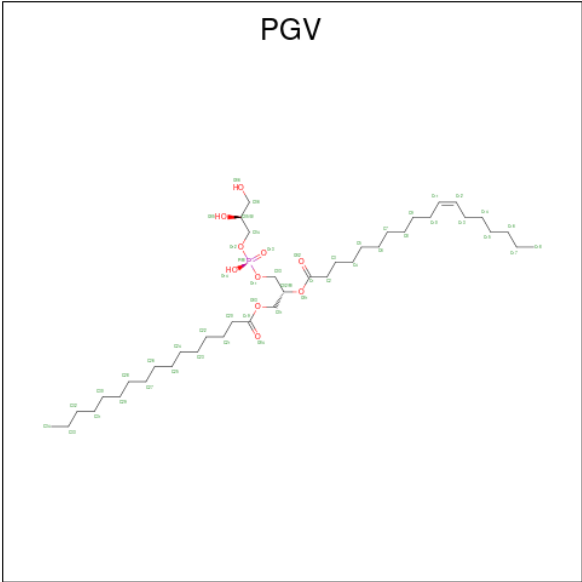
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
18	A	1	Total	C	O	0	0
			2	1	1		
18	N	1	Total	C	O	0	0
			2	1	1		

- Molecule 19 is TRISTEAROYLGLYCEROL (CCD ID: TGL) (formula: C₅₇H₁₁₀O₆).



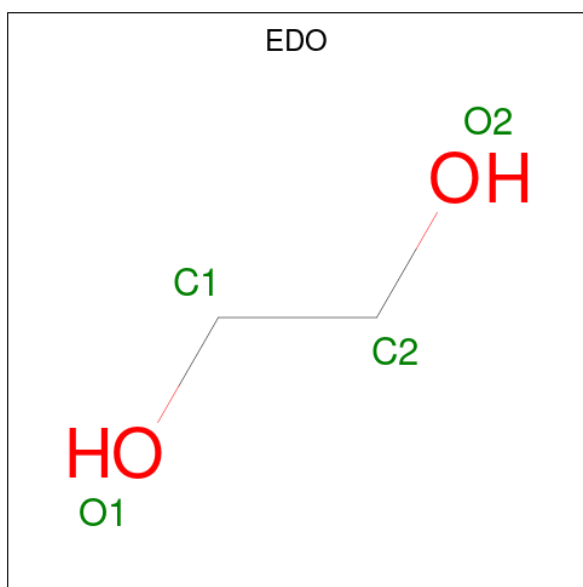
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
19	A	1	Total	C	O	0	0
			63	57	6		
19	D	1	Total	C	O	0	0
			63	57	6		
19	L	1	Total	C	O	0	0
			63	57	6		
19	O	1	Total	C	O	0	0
			63	57	6		
19	Q	1	Total	C	O	0	0
			63	57	6		
19	Y	1	Total	C	O	0	0
			63	57	6		

- Molecule 20 is (1R)-2-{{{[(2S)-2,3-DIHYDROXYPROPYL]OXY}(HYDROXY)PHOSPHORYL]OXY}-1-[(PALMITOYLOXY)METHYL]ETHYL (11E)-OCTADEC-11-ENOATE (CCD ID: PGV) (formula: C₄₀H₇₇O₁₀P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
20	A	1	Total	C	O	P	0	0
			51	40	10	1		
20	A	1	Total	C	O	P	0	0
			51	40	10	1		
20	C	1	Total	C	O	P	0	0
			51	40	10	1		
20	C	1	Total	C	O	P	0	0
			51	40	10	1		
20	N	1	Total	C	O	P	0	0
			51	40	10	1		
20	N	1	Total	C	O	P	0	0
			51	40	10	1		
20	P	1	Total	C	O	P	0	0
			51	40	10	1		
20	P	1	Total	C	O	P	0	0
			51	40	10	1		

- Molecule 21 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



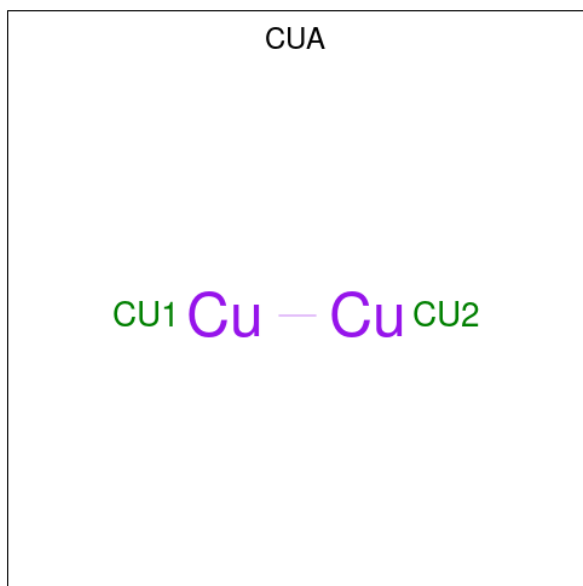
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	A	1	Total	C	O	0	0
			4	2	2		
21	B	1	Total	C	O	0	0
			4	2	2		
21	B	1	Total	C	O	0	0
			4	2	2		
21	C	1	Total	C	O	0	0
			4	2	2		
21	C	1	Total	C	O	0	0
			4	2	2		
21	C	1	Total	C	O	0	0
			4	2	2		
21	C	1	Total	C	O	0	0
			4	2	2		
21	F	1	Total	C	O	0	0
			4	2	2		
21	F	1	Total	C	O	0	0
			4	2	2		
21	G	1	Total	C	O	0	0
			4	2	2		
21	H	1	Total	C	O	0	0
			4	2	2		

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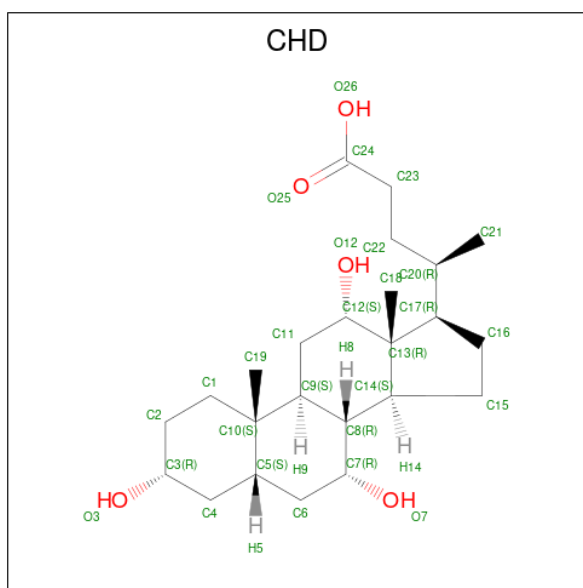
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
21	K	1	Total C O 4 2 2	0	0
21	K	1	Total C O 4 2 2	0	0
21	L	1	Total C O 4 2 2	0	0
21	N	1	Total C O 4 2 2	0	0
21	N	1	Total C O 4 2 2	0	0
21	O	1	Total C O 4 2 2	0	0
21	P	1	Total C O 4 2 2	0	0
21	P	1	Total C O 4 2 2	0	0
21	T	1	Total C O 4 2 2	0	0

- Molecule 22 is DINUCLEAR COPPER ION (CCD ID: CUA) (formula: Cu₂).



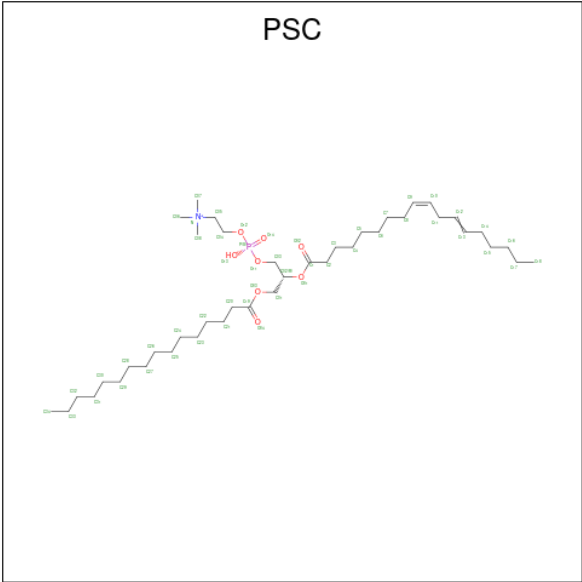
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
22	B	1	Total Cu 2 2	0	0
22	O	1	Total Cu 2 2	0	0

- Molecule 23 is CHOLIC ACID (CCD ID: CHD) (formula: $C_{24}H_{40}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
23	B	1	Total	C	O	0	0
			29	24	5		
23	C	1	Total	C	O	0	0
			29	24	5		
23	C	1	Total	C	O	0	0
			29	24	5		
23	G	1	Total	C	O	0	0
			29	24	5		
23	J	1	Total	C	O	0	0
			29	24	5		
23	P	1	Total	C	O	0	0
			29	24	5		
23	P	1	Total	C	O	0	0
			29	24	5		
23	W	1	Total	C	O	0	0
			29	24	5		

- Molecule 24 is (7R,17E,20E)-4-HYDROXY-N,N,N-TRIMETHYL-9-OXO-7-[(PALMITOYLOXY)METHYL]-3,5,8-TRIOXA-4-PHOSPHAHEXACOSA-17,20-DIEN-1-AMINIUM 4-OXIDE (CCD ID: PSC) (formula: $C_{42}H_{81}NO_8P$).

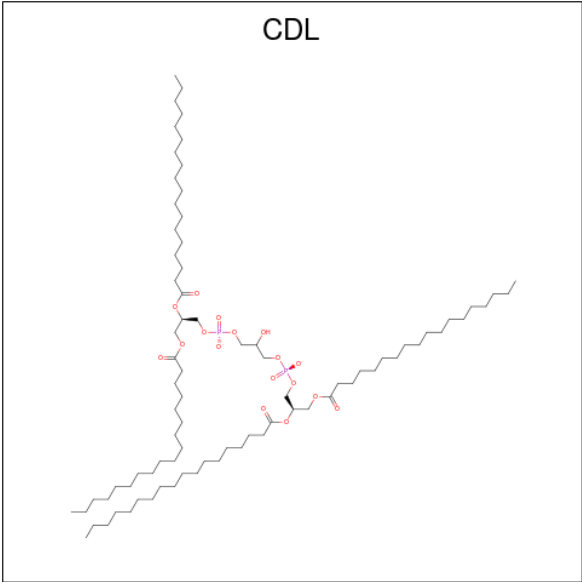


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
24	B	1	Total	C	N	O	P	0	0
			52	42	1	8	1		
24	O	1	Total	C	N	O	P	0	0
			52	42	1	8	1		

- Molecule 25 is UNKNOWN ATOM OR ION (CCD ID: UNX) (formula: X).

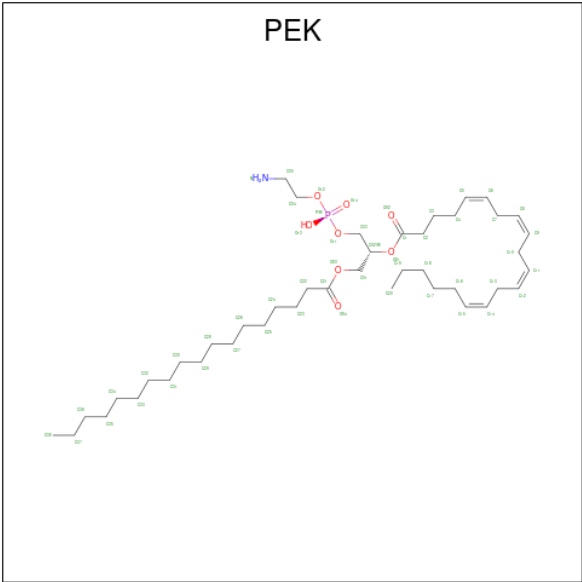
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
25	C	1	Total	X	0	0
			1	1		
25	P	1	Total	X	0	0
			1	1		

- Molecule 26 is CARDIOLIPIN (CCD ID: CDL) (formula: C₈₁H₁₅₆O₁₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
26	C	1	Total	C	O	P	0	0
			100	81	17	2		
26	G	1	Total	C	O	P	0	0
			100	81	17	2		
26	P	1	Total	C	O	P	0	0
			100	81	17	2		
26	T	1	Total	C	O	P	0	0
			100	81	17	2		

- Molecule 27 is (1S)-2-{[(2-AMINOETHOXY)(HYDROXY)PHOSPHORYL]OXY}-1-[(STEAROYLOXY)METHYL]ETHYL (5E,8E,11E,14E)-ICOSA-5,8,11,14-TETRAENOATE (CCD ID: PEK) (formula: C₄₃H₇₈NO₈P).

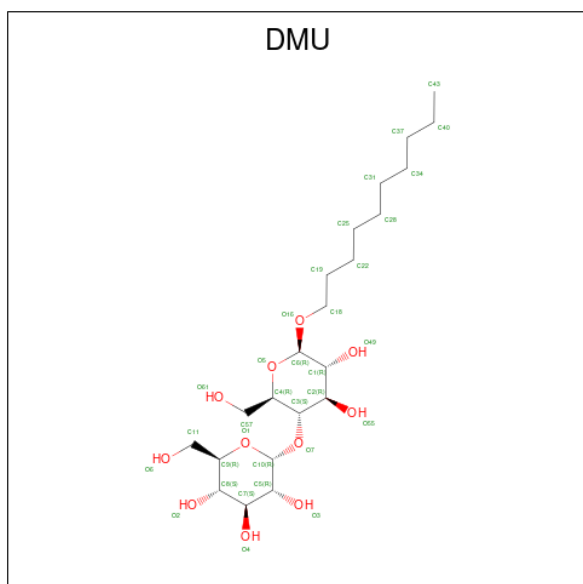


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
27	C	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
27	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
27	G	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
27	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
27	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		
27	T	1	Total	C	N	O	P	0	0
			53	43	1	8	1		

- Molecule 28 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
28	F	1	Total	Zn	0	0
			1	1		
28	S	1	Total	Zn	0	0
			1	1		

- Molecule 29 is DECYL-BETA-D-MALTOPYRANOSIDE (CCD ID: DMU) (formula: C₂₂H₄₂O₁₁).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
29	M	1	Total	C	O	0	0
			33	22	11		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
29	Z	1	Total	C	O	0	0
			33	22	11		

- Molecule 30 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
30	A	191	Total	O	0	0
			191	191		
30	B	120	Total	O	0	1
			121	121		
30	C	83	Total	O	0	0
			83	83		
30	D	89	Total	O	0	0
			89	89		
30	E	45	Total	O	0	0
			45	45		
30	F	49	Total	O	0	0
			49	49		
30	G	40	Total	O	0	0
			40	40		
30	H	37	Total	O	0	0
			37	37		
30	I	19	Total	O	0	0
			19	19		
30	J	19	Total	O	0	0
			19	19		
30	K	27	Total	O	0	0
			27	27		
30	L	27	Total	O	0	0
			27	27		
30	M	18	Total	O	0	0
			18	18		
30	N	169	Total	O	0	0
			169	169		
30	O	100	Total	O	0	1
			101	101		
30	P	91	Total	O	0	0
			91	91		
30	Q	38	Total	O	0	0
			38	38		
30	R	40	Total	O	0	0
			40	40		

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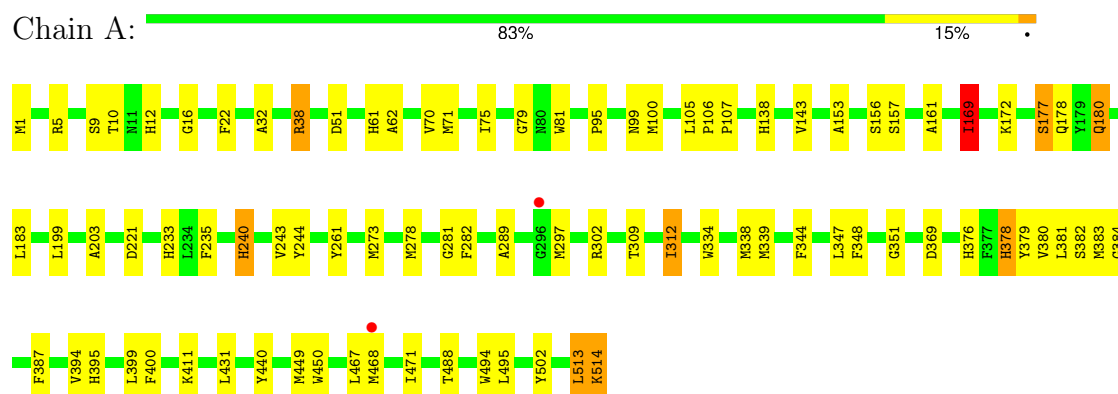
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
30	S	55	Total 55	O 55	0	0
30	T	45	Total 45	O 45	0	0
30	U	27	Total 27	O 27	0	0
30	V	23	Total 23	O 23	0	0
30	W	23	Total 23	O 23	0	0
30	X	16	Total 16	O 16	0	0
30	Y	18	Total 18	O 18	0	0
30	Z	4	Total 4	O 4	0	0

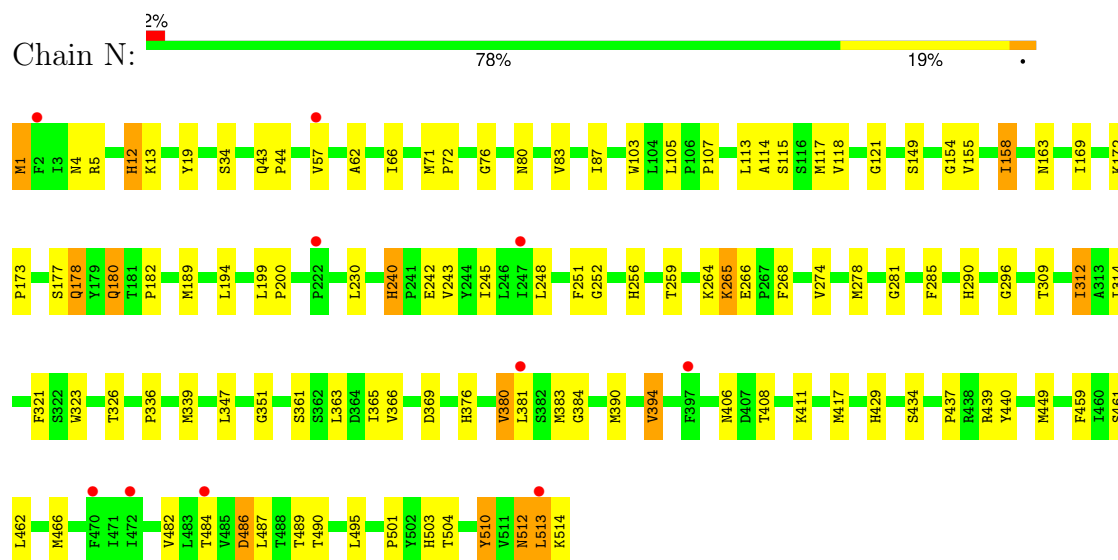
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

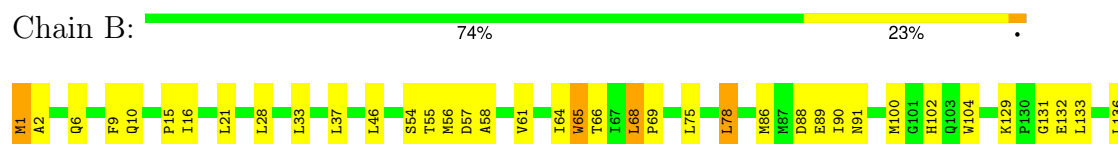
• Molecule 1: Cytochrome c oxidase subunit 1



• Molecule 1: Cytochrome c oxidase subunit 1

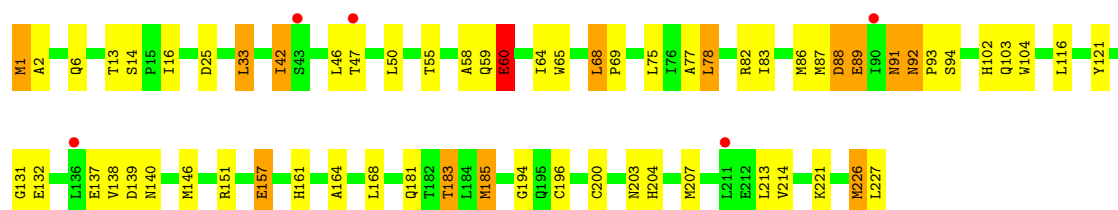


• Molecule 2: Cytochrome c oxidase subunit 2

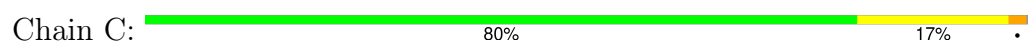




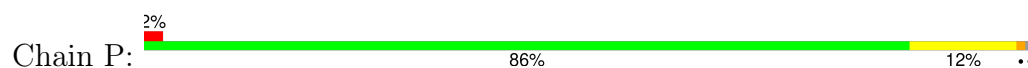
• Molecule 2: Cytochrome c oxidase subunit 2



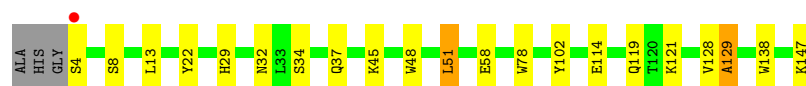
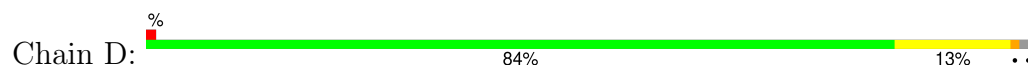
• Molecule 3: Cytochrome c oxidase subunit 3



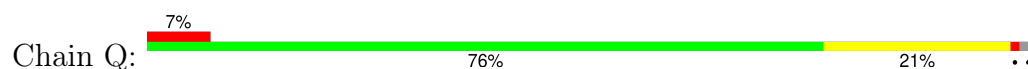
• Molecule 3: Cytochrome c oxidase subunit 3

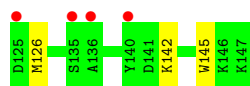


• Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial

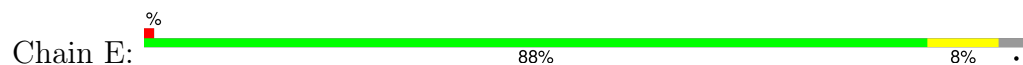


• Molecule 4: Cytochrome c oxidase subunit 4 isoform 1, mitochondrial

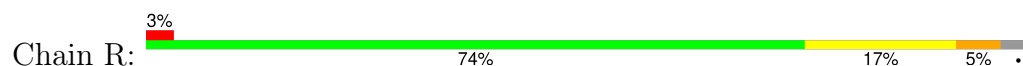




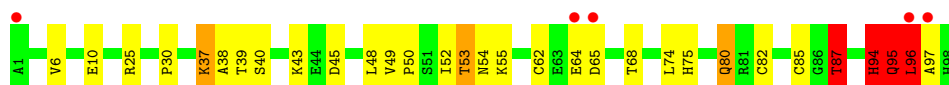
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



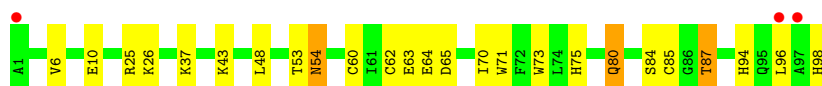
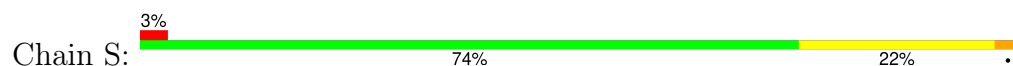
- Molecule 5: Cytochrome c oxidase subunit 5A, mitochondrial



- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



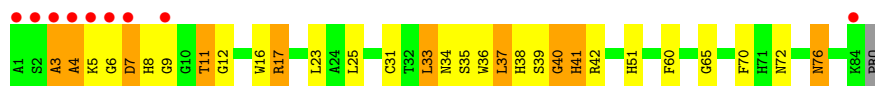
- Molecule 6: Cytochrome c oxidase subunit 5B, mitochondrial



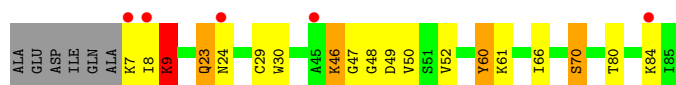
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



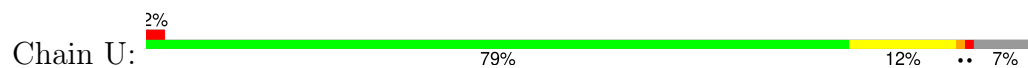
- Molecule 7: Cytochrome c oxidase subunit 6A2, mitochondrial



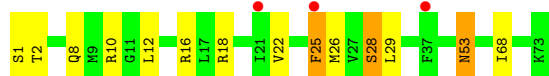
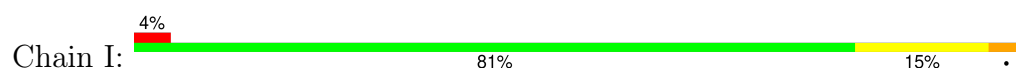
- Molecule 8: Cytochrome c oxidase subunit 6B1



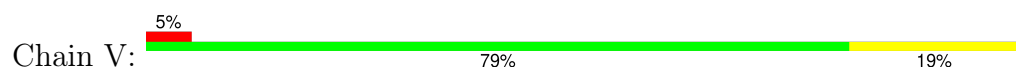
- Molecule 8: Cytochrome c oxidase subunit 6B1



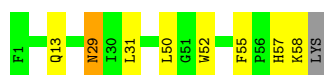
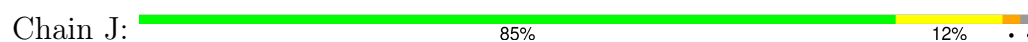
- Molecule 9: Cytochrome c oxidase subunit 6C



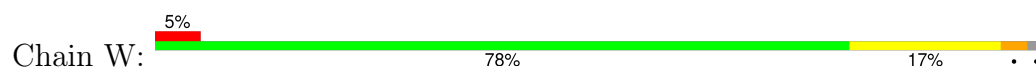
- Molecule 9: Cytochrome c oxidase subunit 6C



- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



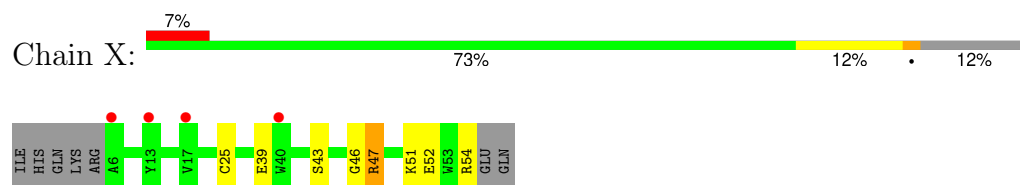
- Molecule 10: Cytochrome c oxidase subunit 7A1, mitochondrial



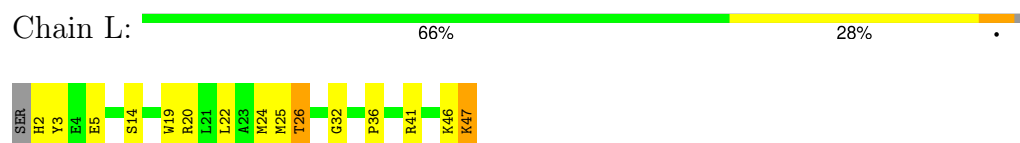
- Molecule 11: Cytochrome c oxidase subunit 7B, mitochondrial



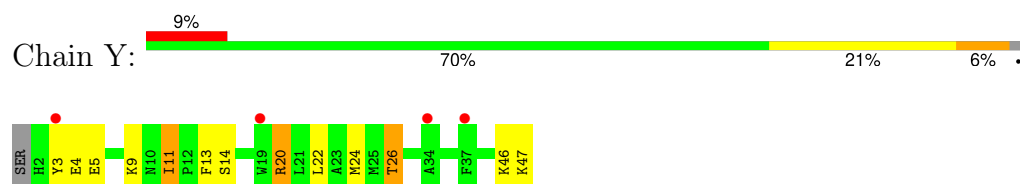
- Molecule 11: Cytochrome c oxidase subunit 7B, mitochondrial



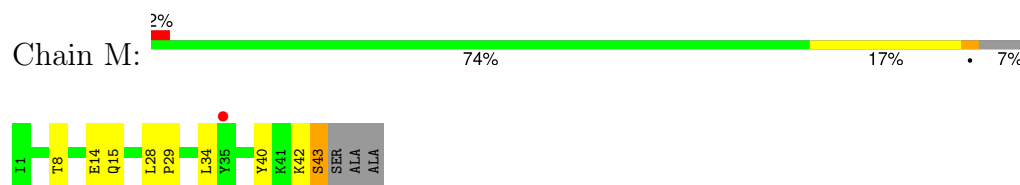
- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



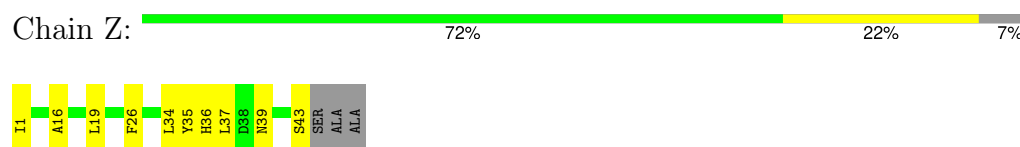
- Molecule 12: Cytochrome c oxidase subunit 7C, mitochondrial



- Molecule 13: Cytochrome c oxidase subunit 8B, mitochondrial



- Molecule 13: Cytochrome c oxidase subunit 8B, mitochondrial



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	185.03Å 208.85Å 178.05Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.20 25.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	96.8 (25.00-2.20) 96.8 (25.00-2.20)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.19Å)	Xtriage
Refinement program	REFMAC 5.8.0048	Depositor
R, R_{free}	0.173 , 0.210 0.174 , 0.210	Depositor DCC
R_{free} test set	17156 reflections (1.36%)	wwPDB-VP
Wilson B-factor (Å ²)	51.4	Xtriage
Anisotropy	0.184	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.21 , 35.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	32242	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.92% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CMO, PSC, CU, MG, UNX, FME, CHD, TGL, PGV, TPO, PEK, NA, HEA, ZN, CDL, DMU, SAC, EDO, CUA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.34	15/4205 (0.4%)	1.23	16/5744 (0.3%)
1	N	1.29	15/4156 (0.4%)	1.22	16/5678 (0.3%)
2	B	1.21	3/1860 (0.2%)	1.25	6/2534 (0.2%)
2	O	1.08	1/1860 (0.1%)	1.17	7/2534 (0.3%)
3	C	1.29	5/2197 (0.2%)	1.17	7/3005 (0.2%)
3	P	1.29	9/2197 (0.4%)	1.14	6/3005 (0.2%)
4	D	1.13	1/1229 (0.1%)	1.13	6/1658 (0.4%)
4	Q	0.93	0/1229	1.06	1/1658 (0.1%)
5	E	1.06	0/871	1.15	0/1182
5	R	0.92	0/871	1.06	0/1182
6	F	1.16	0/765	1.17	2/1038 (0.2%)
6	S	1.12	0/765	1.14	2/1038 (0.2%)
7	G	1.32	1/690 (0.1%)	1.25	6/937 (0.6%)
7	T	1.23	2/690 (0.3%)	1.13	1/937 (0.1%)
8	H	1.06	1/682 (0.1%)	1.16	0/921
8	U	0.96	0/682	1.07	1/921 (0.1%)
9	I	0.94	0/617	1.16	0/818
9	V	0.88	0/629	1.05	0/834
10	J	1.09	0/471	1.17	1/636 (0.2%)
10	W	0.93	0/471	0.99	0/636
11	K	1.17	1/398 (0.3%)	1.10	2/546 (0.4%)
11	X	1.01	0/398	1.15	1/546 (0.2%)
12	L	1.18	0/393	1.18	0/526
12	Y	1.04	0/393	1.08	1/526 (0.2%)
13	M	1.09	0/345	1.15	2/470 (0.4%)
13	Z	0.94	0/345	1.07	0/470
All	All	1.18	54/29409 (0.2%)	1.17	84/39980 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected

by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
8	H	0	1

All (54) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	376	HIS	CE1-NE2	8.38	1.41	1.32
1	A	376	HIS	ND1-CE1	8.01	1.40	1.32
1	A	376	HIS	CG-CD2	7.72	1.44	1.35
1	A	156	SER	CA-C	6.87	1.61	1.52
2	B	204	HIS	ND1-CE1	6.44	1.39	1.32
1	A	61	HIS	ND1-CE1	6.24	1.38	1.32
3	C	14	SER	C-O	-6.14	1.15	1.25
3	P	231	HIS	C-O	6.11	1.31	1.24
1	A	273	MET	N-CA	6.07	1.53	1.46
1	N	149	SER	CA-C	6.03	1.60	1.52
3	P	204	HIS	CG-CD2	5.95	1.42	1.35
1	A	12	HIS	ND1-CE1	5.90	1.38	1.32
1	N	376	HIS	CE1-NE2	5.81	1.38	1.32
1	N	503	HIS	CA-C	-5.79	1.45	1.52
1	A	395	HIS	CG-CD2	5.78	1.42	1.35
3	P	232	HIS	CG-CD2	5.76	1.42	1.35
1	N	251	PHE	C-O	5.72	1.30	1.24
1	A	378	HIS	ND1-CE1	5.72	1.38	1.32
1	N	12	HIS	CE1-NE2	5.66	1.38	1.32
11	K	41	ASN	CA-C	5.64	1.57	1.53
1	N	512	ASN	CA-C	-5.58	1.45	1.52
1	N	256	HIS	CG-CD2	5.55	1.42	1.35
1	A	395	HIS	ND1-CE1	5.49	1.38	1.32
3	C	71	HIS	CG-CD2	5.46	1.41	1.35
1	A	81	TRP	CD2-CE2	5.41	1.50	1.41
1	N	182	PRO	C-O	5.40	1.30	1.23
1	N	163	ASN	CA-C	5.39	1.59	1.52
1	A	235	PHE	CA-C	5.38	1.59	1.52
1	N	290	HIS	CG-CD2	5.38	1.41	1.35
7	T	76	ASN	C-O	-5.34	1.20	1.25
3	P	150	SER	N-CA	5.30	1.52	1.46
3	P	232	HIS	CE1-NE2	5.26	1.37	1.32
7	G	51	HIS	CG-CD2	5.26	1.41	1.35
3	P	103	HIS	CE1-NE2	5.25	1.37	1.32
2	B	65	TRP	CD2-CE2	5.23	1.50	1.41
2	B	102	HIS	CE1-NE2	5.21	1.37	1.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	H	24	ASN	N-CA	-5.20	1.40	1.46
1	N	323	TRP	CD2-CE2	5.19	1.50	1.41
3	P	70	HIS	CE1-NE2	5.19	1.37	1.32
3	P	103	HIS	CG-CD2	5.16	1.41	1.35
1	A	378	HIS	CG-ND1	-5.15	1.32	1.38
1	N	256	HIS	ND1-CE1	5.13	1.37	1.32
2	O	68	LEU	CA-C	5.13	1.58	1.52
1	N	351	GLY	N-CA	5.12	1.52	1.45
3	C	103	HIS	CG-CD2	5.11	1.41	1.35
3	P	91	VAL	N-CA	5.07	1.52	1.46
3	C	232	HIS	CG-CD2	5.06	1.41	1.35
1	A	71	MET	C-O	-5.06	1.19	1.24
1	N	381	LEU	N-CA	5.04	1.52	1.46
1	N	429	HIS	CG-CD2	5.03	1.41	1.35
1	A	16	GLY	N-CA	5.02	1.52	1.45
3	C	240	TRP	CD2-CE2	5.02	1.49	1.41
7	T	51	HIS	CG-CD2	5.02	1.41	1.35
4	D	29	HIS	CE1-NE2	5.00	1.37	1.32

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	70	VAL	N-CA-C	10.20	120.20	110.30
7	T	3	ALA	N-CA-C	8.95	122.47	109.14
1	N	513	LEU	N-CA-C	8.06	119.97	111.03
1	N	105	LEU	CA-C-N	-7.49	112.66	120.38
1	N	105	LEU	C-N-CA	-7.49	112.66	120.38
3	C	129	VAL	CA-C-N	-7.30	111.25	119.28
3	C	129	VAL	C-N-CA	-7.30	111.25	119.28
1	A	169	ILE	CB-CA-C	-7.20	102.29	112.22
2	B	65	TRP	CB-CA-C	6.95	122.73	110.37
3	C	214	PHE	N-CA-C	-6.81	103.94	111.36
2	B	58	ALA	N-CA-C	6.68	120.65	112.23
1	N	83	VAL	N-CA-CB	6.62	114.95	110.52
1	N	248	LEU	CA-C-N	-6.51	112.12	119.28
1	N	248	LEU	C-N-CA	-6.51	112.12	119.28
7	G	58	LYS	CA-C-N	-6.48	113.11	119.78
7	G	58	LYS	C-N-CA	-6.48	113.11	119.78
1	N	240	HIS	N-CA-CB	6.40	117.98	110.42
11	X	43	SER	N-CA-C	6.37	117.41	109.57
2	B	64	ILE	N-CA-C	-6.36	104.13	110.62
2	B	64	ILE	CA-C-N	-6.34	111.65	122.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	64	ILE	C-N-CA	-6.34	111.65	122.56
1	A	199	LEU	CA-C-N	-6.34	112.47	119.19
1	A	199	LEU	C-N-CA	-6.34	112.47	119.19
7	G	72	ASN	CA-C-N	6.30	126.78	119.47
7	G	72	ASN	C-N-CA	6.30	126.78	119.47
7	G	8	HIS	N-CA-C	6.24	120.16	112.54
1	N	380	VAL	CB-CA-C	6.23	122.32	112.16
12	Y	20	ARG	N-CA-C	-6.15	104.49	111.07
3	P	90	GLU	N-CA-C	-6.11	104.17	111.69
4	D	22	TYR	CA-C-N	6.09	125.98	119.28
4	D	22	TYR	C-N-CA	6.09	125.98	119.28
4	D	102	TYR	CA-C-N	-6.05	116.97	123.08
4	D	102	TYR	C-N-CA	-6.05	116.97	123.08
2	O	226	MET	N-CA-C	-5.94	105.34	113.30
1	N	449	MET	N-CA-C	5.84	117.32	111.07
1	A	71	MET	CG-SD-CE	-5.83	88.08	100.90
2	O	65	TRP	N-CA-C	5.77	119.94	113.02
3	P	39	SER	N-CA-C	5.77	118.90	108.69
7	G	7	ASP	N-CA-C	5.75	123.05	110.80
3	C	125	ASN	CA-C-N	-5.69	113.76	119.56
3	C	125	ASN	C-N-CA	-5.69	113.76	119.56
1	A	240	HIS	CA-CB-CG	-5.67	108.12	113.80
1	A	71	MET	CA-C-N	-5.66	113.19	119.19
1	A	71	MET	C-N-CA	-5.66	113.19	119.19
3	P	113	GLY	N-CA-C	-5.66	107.45	115.43
2	O	58	ALA	N-CA-C	5.63	119.32	112.23
1	A	240	HIS	N-CA-CB	5.59	117.02	110.42
3	P	109	THR	N-CA-C	5.57	118.23	110.20
1	N	169	ILE	N-CA-C	5.57	116.34	110.72
1	A	233	HIS	N-CA-C	5.50	116.95	111.07
2	B	177	GLY	N-CA-C	-5.46	108.10	115.36
4	Q	32	ASN	N-CA-C	5.46	117.57	108.02
1	A	161	ALA	N-CA-C	-5.41	105.46	111.36
1	A	244	TYR	CA-CB-CG	-5.30	104.37	113.90
1	N	240	HIS	CA-CB-CG	-5.29	108.51	113.80
8	U	8	ILE	N-CA-C	5.29	120.33	109.34
3	P	33	MET	N-CA-C	-5.28	105.60	111.36
2	O	214	VAL	CA-C-N	5.22	125.21	119.89
2	O	214	VAL	C-N-CA	5.22	125.21	119.89
1	N	266	GLU	N-CA-C	5.21	114.30	108.25
6	F	82	CYS	N-CA-C	-5.21	103.02	109.64
2	O	64	ILE	CA-C-N	-5.18	113.66	122.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	O	64	ILE	C-N-CA	-5.18	113.66	122.56
11	K	6	ALA	CA-C-N	5.17	126.31	119.84
11	K	6	ALA	C-N-CA	5.17	126.31	119.84
10	J	13	GLN	N-CA-C	-5.15	106.68	113.17
1	N	230	LEU	N-CA-C	-5.15	105.67	111.28
3	C	29	SER	CB-CA-C	-5.11	102.64	110.81
1	A	221	ASP	CA-C-N	5.10	125.39	119.47
1	A	221	ASP	C-N-CA	5.10	125.39	119.47
13	M	8	THR	CA-C-N	5.10	125.10	119.90
13	M	8	THR	C-N-CA	5.10	125.10	119.90
1	N	252	GLY	N-CA-C	-5.09	106.59	112.50
1	A	99	ASN	N-CA-C	-5.07	105.94	111.82
3	C	58	TRP	N-CA-C	5.05	116.87	111.36
6	S	6	VAL	CA-C-N	-5.04	115.04	120.03
6	S	6	VAL	C-N-CA	-5.04	115.04	120.03
3	P	228	THR	N-CA-C	-5.04	103.00	110.46
1	N	376	HIS	CA-C-N	-5.03	112.34	121.14
1	N	376	HIS	C-N-CA	-5.03	112.34	121.14
1	A	281	GLY	N-CA-C	-5.03	106.67	112.50
6	F	87	THR	N-CA-CB	5.02	117.41	109.83
4	D	129	ALA	CA-C-N	5.00	124.61	119.56
4	D	129	ALA	C-N-CA	5.00	124.61	119.56

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
8	H	9	LYS	Peptide

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4068	0	4048	91	0
1	N	4027	0	4001	94	0
2	B	1824	0	1833	38	0
2	O	1824	0	1833	49	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	C	2110	0	2027	46	0
3	P	2110	0	2027	18	0
4	D	1195	0	1183	12	0
4	Q	1195	0	1183	30	0
5	E	852	0	845	5	0
5	R	852	0	845	14	0
6	F	748	0	728	25	0
6	S	748	0	728	17	0
7	G	675	0	643	23	0
7	T	675	0	643	30	0
8	H	662	0	623	11	0
8	U	662	0	623	6	0
9	I	609	0	622	11	0
9	V	617	0	631	7	0
10	J	460	0	459	3	0
10	W	460	0	459	13	0
11	K	384	0	366	6	0
11	X	384	0	366	7	0
12	L	380	0	380	20	0
12	Y	380	0	380	11	0
13	M	335	0	352	7	0
13	Z	335	0	352	8	0
14	A	120	0	108	14	0
14	N	120	0	108	12	0
15	A	1	0	0	0	0
15	N	1	0	0	0	0
16	A	1	0	0	0	0
16	N	1	0	0	0	0
17	A	1	0	0	0	0
17	N	1	0	0	0	0
18	A	2	0	0	1	0
18	N	2	0	0	2	0
19	A	63	0	110	3	0
19	D	63	0	110	5	0
19	L	63	0	110	9	0
19	O	63	0	110	7	0
19	Q	63	0	110	5	0
19	Y	63	0	110	5	0
20	A	102	0	152	5	0
20	C	102	0	152	4	0
20	N	102	0	152	6	0
20	P	102	0	152	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
21	A	16	0	24	18	0
21	B	8	0	12	1	0
21	C	16	0	24	2	0
21	F	8	0	12	2	0
21	G	4	0	6	0	0
21	H	4	0	6	2	0
21	K	8	0	12	0	0
21	L	4	0	6	0	0
21	N	8	0	12	3	0
21	O	4	0	6	0	0
21	P	8	0	12	0	0
21	T	4	0	6	3	0
22	B	2	0	0	0	0
22	O	2	0	0	0	0
23	B	29	0	39	1	0
23	C	58	0	78	5	0
23	G	29	0	39	1	0
23	J	29	0	38	0	0
23	P	58	0	78	6	0
23	W	29	0	38	9	0
24	B	52	0	80	4	0
24	O	52	0	80	5	0
25	C	1	0	0	0	0
25	P	1	0	0	0	0
26	C	100	0	156	6	0
26	G	100	0	156	6	0
26	P	100	0	156	7	0
26	T	100	0	156	21	0
27	C	53	0	77	3	0
27	G	106	0	154	6	0
27	T	159	0	231	13	0
28	F	1	0	0	0	0
28	S	1	0	0	0	0
29	M	33	0	42	0	0
29	Z	33	0	42	2	0
30	A	191	0	0	17	0
30	B	121	0	0	5	0
30	C	83	0	0	16	0
30	D	89	0	0	1	0
30	E	45	0	0	2	0
30	F	49	0	0	3	0
30	G	40	0	0	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
30	H	37	0	0	1	0
30	I	19	0	0	0	0
30	J	19	0	0	1	0
30	K	27	0	0	1	0
30	L	27	0	0	6	0
30	M	18	0	0	0	0
30	N	169	0	0	23	0
30	O	101	0	0	20	0
30	P	91	0	0	7	0
30	Q	38	0	0	8	0
30	R	40	0	0	0	0
30	S	55	0	0	5	0
30	T	45	0	0	3	0
30	U	27	0	0	0	0
30	V	23	0	0	1	0
30	W	23	0	0	4	0
30	X	16	0	0	1	0
30	Y	18	0	0	3	0
30	Z	4	0	0	0	0
All	All	32242	0	31442	640	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 10.

All (640) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:22:PHE:HE1	30:A:867:HOH:O	1.15	1.25
1:N:312:ILE:HB	30:N:844:HOH:O	1.31	1.23
7:G:50:TYR:HA	30:G:208:HOH:O	1.07	1.23
30:A:704:HOH:O	3:C:17:PRO:HB3	1.34	1.21
3:P:246:ASP:HB2	30:P:464:HOH:O	1.35	1.20
1:N:259:THR:HA	30:N:740:HOH:O	1.39	1.20
3:C:160:LEU:HB3	30:C:459:HOH:O	1.43	1.16
3:C:99:TRP:CD1	30:C:408:HOH:O	1.98	1.15
12:L:24:MET:HG3	30:L:224:HOH:O	1.47	1.14
21:A:613:EDO:H21	30:A:762:HOH:O	0.97	1.13
2:O:69:PRO:HB2	30:O:455:HOH:O	1.49	1.13
3:C:21:ALA:HB2	30:C:448:HOH:O	1.49	1.11
1:N:434:SER:HB3	30:O:409:HOH:O	1.51	1.10
2:O:161:HIS:HE1	30:O:478:HOH:O	1.31	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:100:MET:HA	30:A:704:HOH:O	1.52	1.09
14:A:602:HEA:HBC1	14:A:602:HEA:HMC1	1.33	1.08
2:O:200:CYS:HB3	30:O:404:HOH:O	1.53	1.08
30:G:208:HOH:O	8:H:80:THR:HB	1.53	1.06
6:S:85:CYS:SG	6:S:87:THR:HG23	1.94	1.06
2:O:161:HIS:CE1	30:O:478:HOH:O	2.06	1.04
4:D:114:GLU:HG3	30:K:209:HOH:O	1.59	1.03
3:C:51:MET:HA	30:C:444:HOH:O	1.61	1.01
1:A:62:ALA:HA	30:A:846:HOH:O	1.58	1.01
30:A:855:HOH:O	19:D:201:TGL:HB21	1.65	0.97
7:T:5:LYS:HD2	27:T:102:PEK:H383	1.42	0.96
9:I:53:ASN:HD22	9:I:53:ASN:H	1.13	0.96
3:C:160:LEU:CB	30:C:459:HOH:O	2.02	0.95
1:A:502:TYR:OH	21:A:612:EDO:H22	1.68	0.94
1:A:347[A]:LEU:HD13	1:A:383[A]:MET:SD	2.09	0.92
3:C:54:MET:HE3	30:C:444:HOH:O	1.69	0.92
30:A:704:HOH:O	3:C:17:PRO:CB	2.02	0.90
1:A:106:PRO:HA	30:A:867:HOH:O	1.72	0.89
1:A:347[B]:LEU:HD22	1:A:383[B]:MET:SD	2.12	0.89
1:A:153:ALA:O	21:A:613:EDO:H11	1.72	0.89
6:F:68:THR:HG23	21:F:103:EDO:H12	1.54	0.88
1:A:347[B]:LEU:HD21	1:A:383[B]:MET:HB3	1.52	0.88
20:C:307:PGV:H211	30:C:408:HOH:O	1.75	0.87
30:C:439:HOH:O	21:H:101:EDO:H11	1.74	0.87
1:N:265:LYS:HD2	1:N:265:LYS:H	1.38	0.87
7:T:37:LEU:O	7:T:38:HIS:HD2	1.57	0.87
6:S:75:HIS:H	6:S:80:GLN:HE22	1.22	0.86
12:Y:46:LYS:HA	30:Y:210:HOH:O	1.73	0.86
1:A:138:HIS:HB2	30:A:710:HOH:O	1.73	0.86
1:A:143:VAL:HG21	30:A:710:HOH:O	1.75	0.86
7:G:72:ASN:H	7:G:76:ASN:HD22	1.23	0.86
1:A:138:HIS:CB	30:A:710:HOH:O	2.24	0.85
19:L:101:TGL:OC1	19:L:101:TGL:HC41	1.74	0.85
8:H:49:ASP:HB3	30:H:218:HOH:O	1.76	0.85
27:T:103:PEK:H381	30:T:235:HOH:O	1.76	0.84
1:N:390:MET:HA	1:N:390:MET:HE3	1.60	0.83
3:C:146:TRP:HE1	21:C:311:EDO:H12	1.43	0.83
1:N:326:THR:HG22	30:N:740:HOH:O	1.76	0.83
6:F:85:CYS:SG	6:F:87:THR:HG23	2.19	0.82
7:G:76:ASN:HD21	27:G:101:PEK:HN2	1.22	0.82
6:S:94:HIS:O	6:S:96:LEU:HD13	1.79	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:Q:9:GLU:O	4:Q:9:GLU:HG2	1.80	0.82
11:K:6:ALA:HB1	11:K:7:PRO:HD2	1.60	0.82
4:Q:109:HIS:HD2	30:Q:314:HOH:O	1.62	0.82
14:N:601:HEA:HBC1	14:N:601:HEA:HMC1	1.61	0.82
4:Q:116:VAL:HG13	30:Q:302:HOH:O	1.79	0.81
5:E:11:PHE:HA	30:E:209:HOH:O	1.81	0.81
1:A:243:VAL:HG21	18:A:606:CMO:C	2.11	0.81
1:N:312:ILE:CB	30:N:844:HOH:O	2.01	0.80
1:A:347[B]:LEU:CD2	1:A:383[B]:MET:HB3	2.10	0.80
2:B:90:ILE:HA	30:B:444:HOH:O	1.80	0.80
1:A:513:LEU:O	1:A:514:LYS:HB2	1.81	0.79
3:C:29:SER:HB3	3:C:42:LEU:HD13	1.65	0.79
2:O:13:THR:HB	2:O:168:LEU:HD23	1.65	0.79
30:N:784:HOH:O	19:Q:201:TGL:HB32	1.82	0.79
3:P:160:LEU:HD13	23:P:305:CHD:H181	1.66	0.78
7:T:72:ASN:H	7:T:76:ASN:HD22	1.32	0.78
12:L:32:GLY:HA3	30:L:201:HOH:O	1.82	0.78
26:T:104:CDL:H151	26:T:104:CDL:OB4	1.83	0.78
10:W:7:GLU:HB3	30:W:207:HOH:O	1.82	0.77
10:W:14:GLU:HG2	30:W:210:HOH:O	1.83	0.77
3:C:99:TRP:HD1	30:C:408:HOH:O	1.49	0.77
1:N:326:THR:CG2	30:N:740:HOH:O	2.29	0.76
26:T:104:CDL:H121	26:T:104:CDL:H371	1.67	0.76
26:T:104:CDL:H273	30:T:235:HOH:O	1.84	0.76
1:A:502:TYR:HH	21:A:612:EDO:H22	1.50	0.76
21:A:612:EDO:H12	3:C:13:PRO:HG3	1.67	0.76
12:Y:20:ARG:HD2	30:Y:205:HOH:O	1.86	0.75
2:B:129:LYS:O	2:B:132:GLU:HB2	1.86	0.75
30:G:207:HOH:O	1:N:178:GLN:CB	2.33	0.75
11:X:47:ARG:HB3	11:X:47:ARG:NH1	2.01	0.75
20:A:608:PGV:H61	3:C:54:MET:HG2	1.68	0.74
1:A:100:MET:CA	30:A:704:HOH:O	2.19	0.74
1:N:314:ILE:HG12	30:O:455:HOH:O	1.86	0.74
9:I:53:ASN:HD22	9:I:53:ASN:N	1.83	0.74
6:S:10:GLU:HB2	30:S:224:HOH:O	1.86	0.74
1:N:383:MET:C	30:N:701:HOH:O	2.29	0.73
23:P:305:CHD:H162	23:P:305:CHD:H232	1.69	0.73
6:F:75:HIS:H	6:F:80:GLN:HE22	1.36	0.73
7:T:12:GLY:HA3	30:T:225:HOH:O	1.87	0.73
9:V:61:GLU:HG3	9:V:65:LYS:NZ	2.04	0.73
4:D:34:SER:H	4:D:37:GLN:HE21	1.37	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:I:53:ASN:H	9:I:53:ASN:ND2	1.86	0.72
8:U:10:ASN:H	8:U:10:ASN:HD22	1.37	0.72
4:Q:109:HIS:CD2	30:Q:314:HOH:O	2.38	0.71
4:D:34:SER:H	4:D:37:GLN:NE2	1.89	0.71
26:T:104:CDL:H161	26:T:104:CDL:H421	1.71	0.71
6:F:64:GLU:O	6:F:65:ASP:HB2	1.89	0.71
12:Y:3:TYR:HA	30:Y:208:HOH:O	1.91	0.70
4:Q:102:TYR:HD1	13:Z:35:TYR:CE2	2.10	0.70
11:X:47:ARG:HB3	11:X:47:ARG:CZ	2.22	0.70
12:Y:14:SER:O	12:Y:20:ARG:NH2	2.25	0.69
30:G:207:HOH:O	1:N:178:GLN:HB3	1.90	0.69
4:Q:102:TYR:CD1	13:Z:35:TYR:HE2	2.09	0.69
1:A:502:TYR:OH	21:A:612:EDO:C2	2.40	0.69
12:Y:46:LYS:O	12:Y:47:LYS:HB2	1.92	0.69
1:A:157:SER:OG	21:A:613:EDO:H22	1.93	0.69
2:O:183:THR:CG2	30:O:408:HOH:O	2.40	0.69
23:W:101:CHD:H3	30:W:209:HOH:O	1.93	0.69
2:O:33:LEU:HD23	9:V:28:SER:HB3	1.75	0.68
1:A:381[B]:LEU:HB2	14:A:602:HEA:CAC	2.24	0.68
6:F:10:GLU:OE1	6:F:25:ARG:NH1	2.26	0.68
6:S:85:CYS:SG	6:S:87:THR:CG2	2.79	0.68
12:L:20:ARG:NH2	19:L:101:TGL:HC32	2.08	0.68
23:W:101:CHD:O12	23:W:101:CHD:H222	1.93	0.68
1:A:10:THR:H	21:A:612:EDO:C2	2.07	0.68
4:Q:34:SER:H	4:Q:37:GLN:HE21	1.40	0.68
20:A:609:PGV:H222	13:M:15:GLN:HE22	1.59	0.67
1:N:487:LEU:HD13	30:N:717:HOH:O	1.93	0.67
9:V:61:GLU:HG3	9:V:65:LYS:HZ1	1.57	0.67
11:K:24:PHE:O	11:K:28:VAL:HG12	1.95	0.67
1:A:157:SER:OG	21:A:613:EDO:C2	2.43	0.67
1:A:278:MET:HE2	21:T:105:EDO:O1	1.93	0.67
1:A:502:TYR:OH	21:A:612:EDO:H11	1.95	0.66
4:Q:98:TRP:CZ3	29:Z:101:DMU:H15	2.31	0.66
2:O:88:ASP:HB3	30:O:469:HOH:O	1.95	0.66
4:Q:119:GLN:HB2	30:Q:305:HOH:O	1.96	0.66
6:F:85:CYS:SG	6:F:87:THR:CG2	2.83	0.65
4:Q:102:TYR:CD1	13:Z:35:TYR:CE2	2.82	0.65
10:J:55:PHE:HB2	30:J:208:HOH:O	1.95	0.65
12:L:20:ARG:HH22	19:L:101:TGL:HC52	1.62	0.65
30:B:431:HOH:O	7:T:17:ARG:HD2	1.96	0.65
5:R:56:ARG:NH2	5:R:59:ASN:OD1	2.27	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:30:PRO:O	6:F:96:LEU:HD23	1.97	0.65
12:Y:24:MET:HE1	19:Y:101:TGL:HC21	1.78	0.65
12:L:46:LYS:O	12:L:47:LYS:HB2	1.97	0.65
20:P:303:PGV:H172	26:P:304:CDL:H611	1.77	0.65
26:C:303:CDL:H522	26:C:303:CDL:HB62	1.78	0.64
1:N:274:VAL:HG12	1:N:278:MET:HE2	1.78	0.64
27:T:101:PEK:H32	27:T:101:PEK:H71	1.79	0.64
20:C:307:PGV:C21	30:C:408:HOH:O	2.40	0.64
30:G:207:HOH:O	1:N:178:GLN:HB2	1.96	0.64
2:O:200:CYS:HB2	30:O:478:HOH:O	1.97	0.64
1:A:100:MET:HG3	30:C:448:HOH:O	1.96	0.64
8:U:10:ASN:H	8:U:10:ASN:ND2	1.95	0.64
21:A:612:EDO:H11	12:L:3:TYR:HE2	1.62	0.64
7:T:37:LEU:O	7:T:38:HIS:CD2	2.45	0.64
23:P:306:CHD:H12	23:P:306:CHD:H212	1.80	0.64
27:T:102:PEK:H381	26:T:104:CDL:C87	2.27	0.63
14:A:601:HEA:CHD	30:A:846:HOH:O	2.46	0.63
26:T:104:CDL:H211	26:T:104:CDL:CB5	2.28	0.63
3:C:111:GLU:HG3	21:H:101:EDO:O1	1.98	0.62
2:O:2:ALA:HA	2:O:6:GLN:OE1	1.98	0.62
2:O:161:HIS:HA	30:O:404:HOH:O	1.99	0.62
7:T:5:LYS:HD2	27:T:102:PEK:C38	2.23	0.62
7:T:76:ASN:HD21	27:T:101:PEK:HN2	1.47	0.62
26:T:104:CDL:H522	26:T:104:CDL:H712	1.82	0.62
14:A:601:HEA:HMC1	14:A:601:HEA:HBC1	1.82	0.62
6:F:95:GLN:HG3	6:F:96:LEU:N	2.15	0.62
3:C:51:MET:HE2	30:C:444:HOH:O	1.99	0.62
24:O:303:PSC:H211	24:O:303:PSC:C01	2.30	0.62
2:B:33:LEU:CD1	9:I:28:SER:HB3	2.29	0.61
27:C:306:PEK:H5	27:C:306:PEK:H221	1.82	0.61
1:N:312:ILE:CG2	30:N:844:HOH:O	2.42	0.61
2:O:91:ASN:ND2	30:O:402:HOH:O	2.33	0.61
1:A:62:ALA:HB2	14:A:601:HEA:HBD1	1.83	0.61
3:C:44:MET:HE2	3:C:44:MET:HA	1.82	0.61
20:A:609:PGV:H222	13:M:15:GLN:NE2	2.14	0.61
14:N:601:HEA:HBC1	14:N:601:HEA:CMC	2.31	0.61
1:A:10:THR:H	21:A:612:EDO:H21	1.66	0.61
1:N:296:GLY:HA2	8:U:23:GLN:NE2	2.15	0.61
8:U:10:ASN:HD22	8:U:10:ASN:N	1.96	0.61
24:O:303:PSC:H231	24:O:303:PSC:H011	1.83	0.60
4:Q:98:TRP:CE3	29:Z:101:DMU:H15	2.36	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:F:95:GLN:O	6:F:97:ALA:N	2.33	0.60
1:A:339:MET:HE1	1:A:411:LYS:HD3	1.82	0.60
4:Q:52:SER:OG	4:Q:55:GLU:HG3	2.00	0.60
1:A:105:LEU:HD21	21:A:613:EDO:H12	1.82	0.60
4:D:78:TRP:HA	19:D:201:TGL:HB42	1.83	0.60
1:N:199:LEU:N	1:N:200:PRO:CD	2.64	0.60
1:A:95:PRO:HG2	3:C:11:VAL:HG23	1.83	0.60
3:C:67:PHE:HE2	26:C:303:CDL:H1	1.66	0.60
2:O:33:LEU:CD2	9:V:28:SER:HB3	2.32	0.60
2:O:59:GLN:C	2:O:60:GLU:HG3	2.27	0.60
2:O:131:GLY:HA2	30:Q:305:HOH:O	2.02	0.59
6:S:10:GLU:CB	30:S:224:HOH:O	2.47	0.59
1:A:9:SER:HA	21:A:612:EDO:H21	1.84	0.59
1:A:177:SER:H	1:A:180:GLN:NE2	2.01	0.59
6:F:94:HIS:O	6:F:96:LEU:HD12	2.02	0.59
11:K:6:ALA:HB1	11:K:7:PRO:CD	2.33	0.59
12:L:22:LEU:O	12:L:26:THR:HB	2.01	0.59
7:T:3:ALA:HB1	27:T:102:PEK:H331	1.83	0.59
2:B:89:GLU:O	2:B:91:ASN:OD1	2.21	0.59
1:N:1:FME:HCN	1:N:4:ASN:H	1.68	0.59
1:A:513:LEU:O	1:A:514:LYS:CB	2.49	0.59
24:B:303:PSC:H062	9:I:10:ARG:HE	1.65	0.59
1:N:383:MET:CA	30:N:701:HOH:O	2.51	0.59
3:C:160:LEU:HD13	23:C:304:CHD:H181	1.85	0.58
1:A:138:HIS:HB3	30:A:710:HOH:O	1.96	0.58
12:L:32:GLY:CA	30:L:201:HOH:O	2.46	0.58
7:T:3:ALA:CB	27:T:102:PEK:H331	2.32	0.58
1:N:406:ASN:HD21	20:N:607:PGV:H22	1.68	0.58
1:N:265:LYS:HD2	1:N:265:LYS:N	2.16	0.58
1:A:383[A]:MET:HA	1:A:387:PHE:CD1	2.39	0.58
4:Q:9:GLU:O	4:Q:9:GLU:CG	2.52	0.58
1:A:95:PRO:HG2	3:C:11:VAL:CG2	2.34	0.57
6:S:54:ASN:HD22	6:S:54:ASN:C	2.12	0.57
7:G:4:ALA:C	7:G:5:LYS:HG3	2.30	0.57
10:J:29:ASN:HD22	10:J:29:ASN:H	1.51	0.57
20:N:607:PGV:H02	20:N:607:PGV:H042	1.85	0.57
1:N:365:ILE:HB	30:N:705:HOH:O	2.04	0.57
1:N:265:LYS:HB2	1:N:490:THR:HG21	1.85	0.57
1:N:177:SER:H	1:N:180:GLN:NE2	2.03	0.57
1:N:103:TRP:O	1:N:107:PRO:HD2	2.05	0.57
1:N:243:VAL:HG21	18:N:606:CMO:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
26:P:304:CDL:H341	26:P:304:CDL:H171	1.87	0.56
3:C:3:HIS:HA	30:C:458:HOH:O	2.04	0.56
9:I:53:ASN:N	9:I:53:ASN:ND2	2.48	0.56
4:Q:78:TRP:CE2	4:Q:79:LYS:HG3	2.40	0.56
7:G:6:GLY:HA3	1:N:189:MET:HE3	1.88	0.56
1:N:514:LYS:NZ	6:S:60:CYS:SG	2.77	0.56
2:B:1:FME:HA	21:B:304:EDO:H22	1.87	0.56
4:Q:12:ALA:HA	6:S:73:TRP:CD1	2.41	0.56
1:A:334:TRP:CH2	2:B:46:LEU:HD13	2.41	0.56
2:O:139:ASP:OD1	2:O:140:ASN:N	2.38	0.56
2:O:82:ARG:NH1	2:O:86:MET:HE1	2.21	0.55
9:I:25:PHE:CD1	9:I:25:PHE:C	2.84	0.55
19:A:607:TGL:H341	19:A:607:TGL:H211	1.88	0.55
3:C:154:GLY:HA2	6:F:6:VAL:HB	1.87	0.55
8:H:60:TYR:CD1	8:H:60:TYR:C	2.84	0.55
24:O:303:PSC:H211	24:O:303:PSC:H011	1.89	0.55
5:R:6:GLU:HB2	5:R:10:GLU:OE2	2.05	0.55
14:N:602:HEA:C1D	18:N:606:CMO:O	2.55	0.55
4:Q:145:TRP:CD1	11:X:46:GLY:HA2	2.41	0.55
26:T:104:CDL:HB22	26:T:104:CDL:H131	1.89	0.55
9:V:22:VAL:O	9:V:26:MET:HG2	2.07	0.55
7:G:31:CYS:SG	26:G:102:CDL:H522	2.47	0.55
5:R:72:LYS:NZ	5:R:102:GLU:OE2	2.39	0.55
3:C:157:LYS:HD3	27:C:306:PEK:H052	1.89	0.55
14:N:602:HEA:HBC1	14:N:602:HEA:HMC1	1.89	0.55
3:C:146:TRP:NE1	21:C:311:EDO:H12	2.19	0.54
7:T:38:HIS:NE2	26:T:104:CDL:H122	2.23	0.54
12:L:20:ARG:HH22	19:L:101:TGL:HC32	1.71	0.54
2:O:196:CYS:HB2	2:O:207:MET:HG3	1.89	0.54
10:W:33:ARG:HG2	23:W:101:CHD:H181	1.88	0.54
1:A:502:TYR:OH	21:A:612:EDO:C1	2.55	0.54
2:O:89:GLU:HB3	30:O:433:HOH:O	2.06	0.54
5:R:37:VAL:HG11	5:R:70:VAL:HG21	1.90	0.54
7:T:38:HIS:CE1	26:T:104:CDL:HA21	2.43	0.54
7:T:37:LEU:C	7:T:38:HIS:HD2	2.15	0.54
3:C:50:ASN:ND2	3:C:54:MET:HE2	2.22	0.54
4:Q:116:VAL:HA	30:Q:306:HOH:O	2.08	0.54
19:O:302:TGL:C28	19:O:302:TGL:HB92	2.38	0.54
12:Y:24:MET:HE1	19:Y:101:TGL:CC2	2.37	0.54
7:G:18:PHE:CE1	1:N:278:MET:HE1	2.43	0.53
12:L:19:TRP:HZ2	13:M:14:GLU:HG2	1.72	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
27:T:102:PEK:H381	26:T:104:CDL:H871	1.88	0.53
1:A:172:LYS:HB2	30:A:788:HOH:O	2.07	0.53
1:A:289:ALA:HB1	1:A:297:MET:HE1	1.90	0.53
14:A:602:HEA:HBC1	14:A:602:HEA:CMC	2.20	0.53
7:G:5:LYS:HB2	27:G:103:PEK:H351	1.90	0.53
5:R:82:TYR:HB3	5:R:83:PRO:HD3	1.91	0.53
1:A:400:PHE:HB3	19:L:101:TGL:H283	1.90	0.53
4:D:34:SER:N	4:D:37:GLN:HE21	2.07	0.53
1:A:334:TRP:HH2	2:B:46:LEU:HD13	1.73	0.53
1:N:19:TYR:CD1	1:N:76:GLY:HA3	2.44	0.53
1:N:177:SER:H	1:N:180:GLN:HE22	1.55	0.53
14:A:602:HEA:HMC1	14:A:602:HEA:CBC	2.19	0.53
1:N:417:MET:HE3	30:N:841:HOH:O	2.09	0.53
3:P:38:ASN:HA	30:P:405:HOH:O	2.08	0.53
20:C:307:PGV:H41	30:P:469:HOH:O	2.09	0.52
7:G:23:LEU:C	7:G:26:PRO:HD2	2.34	0.52
2:O:102:HIS:HE1	2:O:157:GLU:OE2	1.92	0.52
4:Q:34:SER:H	4:Q:37:GLN:NE2	2.05	0.52
10:W:33:ARG:CZ	23:W:101:CHD:H162	2.39	0.52
2:O:47:THR:HB	19:Q:201:TGL:H182	1.91	0.52
2:O:103:GLN:HG2	30:O:478:HOH:O	2.10	0.52
2:B:151:ARG:HD3	2:B:181:GLN:HE21	1.74	0.52
2:B:151:ARG:CD	2:B:181:GLN:HE21	2.23	0.52
3:C:47:LEU:O	3:C:51:MET:HG2	2.10	0.52
26:G:102:CDL:H712	26:G:102:CDL:H512	1.91	0.52
2:O:200:CYS:CB	30:O:404:HOH:O	2.31	0.52
1:N:513:LEU:O	1:N:514:LYS:HB2	2.09	0.52
2:O:151:ARG:HE	2:O:181:GLN:NE2	2.08	0.52
26:T:104:CDL:H522	26:T:104:CDL:H231	1.91	0.52
21:A:612:EDO:H11	12:L:3:TYR:CE2	2.45	0.52
2:B:136:LEU:HB3	2:B:193:TYR:CD2	2.44	0.52
10:J:52:TRP:O	10:J:57:HIS:HE1	1.92	0.52
12:L:19:TRP:CZ2	13:M:14:GLU:HG2	2.45	0.52
19:D:201:TGL:H362	9:I:16:ARG:HE	1.76	0.51
1:A:378:HIS:O	1:A:382[B]:SER:HB2	2.10	0.51
23:C:305:CHD:H152	20:C:307:PGV:H102	1.93	0.51
1:N:155:VAL:HG21	20:N:608:PGV:H142	1.91	0.51
1:N:384:GLY:N	30:N:701:HOH:O	2.42	0.51
1:N:408:THR:HB	20:N:607:PGV:H21	1.92	0.51
1:A:32:ALA:HB3	12:L:36:PRO:HG2	1.91	0.51
20:A:609:PGV:H221	20:A:609:PGV:H012	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:43:GLN:HB2	1:N:44:PRO:HD2	1.92	0.51
5:R:25:ASP:OD1	5:R:28:GLU:HG3	2.11	0.51
1:A:22:PHE:HA	19:L:101:TGL:HB71	1.92	0.51
1:N:484:THR:HG22	21:N:610:EDO:C1	2.41	0.51
2:O:91:ASN:N	30:O:403:HOH:O	2.43	0.51
1:N:484:THR:HG22	21:N:610:EDO:O1	2.11	0.51
2:O:69:PRO:HG2	30:O:416:HOH:O	2.11	0.51
5:R:99:SER:HB2	5:R:104:LEU:HD21	1.93	0.51
1:A:381[B]:LEU:HB2	14:A:602:HEA:HAC	1.93	0.51
26:P:304:CDL:HA4	26:P:304:CDL:H121	1.92	0.51
2:O:132:GLU:HB3	2:O:137:GLU:HG3	1.93	0.50
2:B:54:SER:HB2	30:B:473:HOH:O	2.11	0.50
4:D:128:VAL:O	4:D:129:ALA:C	2.53	0.50
5:R:80:GLU:H	5:R:80:GLU:CD	2.19	0.50
26:T:104:CDL:H111	26:T:104:CDL:HA22	1.93	0.50
23:B:302:CHD:H212	23:B:302:CHD:H12	1.93	0.50
6:F:95:GLN:C	6:F:97:ALA:H	2.19	0.50
3:C:226:HIS:CE1	26:C:303:CDL:HB32	2.46	0.50
7:G:17:ARG:HD2	30:G:220:HOH:O	2.10	0.50
12:L:24:MET:HE3	30:L:222:HOH:O	2.12	0.50
2:O:1:FME:HE3	4:Q:123:MET:HE3	1.93	0.50
5:R:105:GLY:O	5:R:108:LYS:HD3	2.12	0.50
6:S:10:GLU:OE2	6:S:25:ARG:NH2	2.45	0.50
20:N:607:PGV:H292	13:Z:16:ALA:HA	1.94	0.50
6:S:64:GLU:O	6:S:65:ASP:HB2	2.11	0.50
7:G:3:ALA:O	7:G:4:ALA:HB2	2.12	0.49
4:Q:44:GLU:O	4:Q:44:GLU:HG2	2.12	0.49
1:A:282:PHE:HA	7:T:4:ALA:CB	2.42	0.49
3:C:160:LEU:HB2	30:C:459:HOH:O	1.88	0.49
1:N:71:MET:HB2	1:N:72:PRO:HD3	1.93	0.49
7:T:38:HIS:HE1	26:T:104:CDL:HA21	1.77	0.49
1:A:282:PHE:HA	7:T:4:ALA:HB3	1.94	0.49
2:B:100:MET:HE3	2:B:157:GLU:HG3	1.95	0.49
23:G:104:CHD:H12	23:G:104:CHD:H212	1.93	0.49
1:N:487:LEU:CD1	30:N:717:HOH:O	2.58	0.49
2:O:131:GLY:CA	30:Q:305:HOH:O	2.58	0.49
4:Q:142:LYS:HD3	30:Q:325:HOH:O	2.13	0.49
6:S:37:LYS:NZ	30:S:201:HOH:O	2.41	0.49
1:A:38:ARG:HD2	14:A:601:HEA:OMA	2.12	0.49
7:G:18:PHE:CZ	1:N:278:MET:CE	2.95	0.49
1:A:240:HIS:CD2	1:A:240:HIS:C	2.91	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
14:A:601:HEA:HBC1	14:A:601:HEA:CMC	2.42	0.49
8:H:46:LYS:O	8:H:48:GLY:N	2.45	0.49
2:O:13:THR:HB	2:O:168:LEU:CD2	2.40	0.49
4:Q:122:ARG:HG2	4:Q:126:MET:HE2	1.94	0.49
1:A:157:SER:OG	21:A:613:EDO:C1	2.61	0.49
14:N:601:HEA:CHD	30:N:759:HOH:O	2.60	0.49
20:N:607:PGV:H02	20:N:607:PGV:C04	2.43	0.49
1:N:437:PRO:HG2	1:N:440:TYR:CZ	2.47	0.49
2:O:60:GLU:HG2	30:O:485:HOH:O	2.12	0.49
3:P:213:THR:HG23	26:P:304:CDL:H771	1.94	0.49
20:P:301:PGV:H011	30:P:472:HOH:O	2.11	0.49
5:R:90:ARG:HB3	5:R:91:PRO:HD3	1.94	0.49
6:F:53:THR:HB	6:F:54:ASN:H	1.49	0.48
1:N:172:LYS:NZ	1:N:178:GLN:HE22	2.09	0.48
3:C:177:GLN:HA	3:C:177:GLN:OE1	2.12	0.48
7:T:37:LEU:C	7:T:38:HIS:CD2	2.91	0.48
2:O:14:SER:HB3	2:O:185:MET:O	2.13	0.48
1:N:242:GLU:HA	1:N:245:ILE:HD12	1.95	0.48
2:O:16:ILE:HD12	2:O:87:MET:HG2	1.94	0.48
7:T:60:PHE:O	7:T:65:GLY:HA2	2.13	0.48
1:A:347[B]:LEU:HD12	1:A:348:PHE:N	2.28	0.48
2:B:104:TRP:CD2	2:B:203:ASN:HB2	2.49	0.48
3:C:51:MET:HB2	26:C:303:CDL:H381	1.95	0.48
12:L:47:LYS:HE2	13:M:43:SER:OG	2.13	0.48
3:C:40:MET:O	3:C:44:MET:HG2	2.14	0.48
26:C:303:CDL:H352	30:C:478:HOH:O	2.14	0.48
9:I:22:VAL:O	9:I:26:MET:HG2	2.14	0.48
5:R:72:LYS:HB2	5:R:82:TYR:CD1	2.48	0.48
26:C:303:CDL:H673	30:C:481:HOH:O	2.14	0.48
1:N:484:THR:HG22	21:N:610:EDO:H12	1.96	0.48
9:V:27:VAL:O	9:V:31[B]:PHE:HD2	1.96	0.48
1:A:172:LYS:HZ2	1:A:178:GLN:HE22	1.61	0.47
1:N:5:ARG:HA	30:N:709:HOH:O	2.14	0.47
20:P:301:PGV:H031	30:P:472:HOH:O	2.14	0.47
2:B:9:PHE:HB2	2:B:21:LEU:HD21	1.96	0.47
23:C:304:CHD:H12	23:C:304:CHD:H212	1.94	0.47
7:G:4:ALA:HB1	1:N:281:GLY:HA3	1.96	0.47
1:A:278:MET:CE	21:T:105:EDO:O1	2.59	0.47
2:B:56:MET:HA	24:B:303:PSC:H232	1.97	0.47
1:N:19:TYR:CG	1:N:76:GLY:HA3	2.49	0.47
1:N:461:SER:HB2	30:N:814:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:72:ASN:N	7:G:76:ASN:HD22	2.01	0.47
1:N:510:TYR:C	1:N:510:TYR:CD1	2.91	0.47
1:N:510:TYR:OH	1:N:512:ASN:ND2	2.43	0.47
2:O:146:MET:HA	2:O:213:LEU:HD12	1.96	0.47
1:A:380[A]:VAL:O	1:A:384[A]:GLY:HA3	2.14	0.47
12:L:41:ARG:HG3	13:M:40:TYR:CE2	2.50	0.47
1:N:117:MET:HB3	10:W:54:SER:OG	2.14	0.47
1:A:172:LYS:NZ	1:A:178:GLN:HE22	2.12	0.47
11:K:31:TYR:CD1	11:K:35:GLN:HG3	2.49	0.47
12:L:14:SER:H	19:L:101:TGL:HC31	1.80	0.47
8:H:9:LYS:HD3	8:H:9:LYS:HA	1.61	0.47
13:M:28:LEU:HB2	13:M:29:PRO:HD3	1.95	0.47
1:N:62:ALA:HB2	14:N:601:HEA:HBD1	1.97	0.47
14:N:602:HEA:H243	2:O:69:PRO:HB3	1.97	0.47
19:A:607:TGL:H222	19:A:607:TGL:C36	2.45	0.47
4:D:78:TRP:HB3	19:D:201:TGL:HB31	1.97	0.47
6:S:62:CYS:HB3	6:S:85:CYS:HB3	1.97	0.47
1:N:154:GLY:O	1:N:158:ILE:HG13	2.15	0.46
23:W:101:CHD:H21	23:W:101:CHD:H9	1.90	0.46
6:F:40:SER:OG	6:F:45:ASP:HB3	2.15	0.46
10:W:33:ARG:NE	23:W:101:CHD:H212	2.30	0.46
1:A:513:LEU:HD23	1:A:513:LEU:HA	1.64	0.46
1:N:336:PRO:HB2	1:N:394:VAL:HG21	1.97	0.46
1:A:347[B]:LEU:CD2	1:A:383[B]:MET:CB	2.90	0.46
3:C:29:SER:HB3	3:C:42:LEU:CD1	2.39	0.46
1:N:312:ILE:HG22	1:N:312:ILE:O	2.15	0.46
1:N:321:PHE:CZ	24:O:303:PSC:H161	2.51	0.46
8:H:66:ILE:O	8:H:70:SER:HB3	2.16	0.46
1:N:43:GLN:HB2	1:N:44:PRO:CD	2.46	0.46
19:O:302:TGL:HB92	19:O:302:TGL:H283	1.97	0.46
6:F:10:GLU:HB2	30:F:229:HOH:O	2.16	0.46
7:T:70:PHE:HB2	27:T:101:PEK:H041	1.98	0.46
2:B:2:ALA:HA	2:B:6:GLN:OE1	2.16	0.46
2:B:16:ILE:HD13	2:B:16:ILE:HA	1.76	0.46
7:T:33:LEU:O	7:T:34:ASN:C	2.58	0.46
1:A:169:ILE:N	1:A:169:ILE:HD13	2.27	0.45
1:N:439:ARG:O	2:O:204:HIS:HD2	1.99	0.45
1:A:351:GLY:HA3	1:A:380[B]:VAL:HB	1.98	0.45
7:G:18:PHE:CE1	1:N:278:MET:CE	2.99	0.45
1:N:268:PHE:CD1	1:N:268:PHE:C	2.94	0.45
24:O:303:PSC:H252	24:O:303:PSC:H21	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:T:7:ASP:O	7:T:9:GLY:N	2.47	0.45
1:A:282:PHE:HE2	26:T:104:CDL:H262	1.81	0.45
1:A:347[B]:LEU:HD13	1:A:379:TYR:O	2.16	0.45
1:N:501:PRO:HA	12:Y:5:GLU:OE2	2.16	0.45
19:O:302:TGL:H281	19:O:302:TGL:CB9	2.46	0.45
6:S:64:GLU:HG3	30:S:243:HOH:O	2.17	0.45
2:B:222:TRP:CE2	2:B:226:MET:HG3	2.52	0.45
1:A:514:LYS:HA	6:F:38:ALA:HB3	1.98	0.45
1:N:107:PRO:HB3	3:P:25:LEU:HB2	1.99	0.45
2:O:25:ASP:HA	30:O:427:HOH:O	2.17	0.45
6:S:25:ARG:HD3	30:S:244:HOH:O	2.15	0.45
11:X:39:GLU:HB3	30:X:106:HOH:O	2.16	0.45
1:A:449:MET:HE2	11:K:40:TRP:HB3	1.98	0.45
2:B:1:FME:CE	2:B:133:LEU:HD13	2.47	0.45
1:N:486:ASP:HB2	30:N:804:HOH:O	2.16	0.45
26:P:304:CDL:H641	26:P:304:CDL:H271	1.99	0.45
1:A:51:ASP:HB2	2:B:202:SER:O	2.17	0.45
7:T:7:ASP:C	7:T:9:GLY:H	2.24	0.45
3:C:42:LEU:HD23	3:C:42:LEU:HA	1.89	0.45
1:N:34:SER:HB2	14:N:601:HEA:C2B	2.47	0.45
1:N:264:LYS:HE2	30:O:413:HOH:O	2.16	0.45
1:N:347:LEU:HD13	1:N:383:MET:SD	2.57	0.45
19:O:302:TGL:HA71	19:O:302:TGL:H202	1.57	0.45
1:A:468:MET:HG3	30:A:873:HOH:O	2.17	0.44
4:D:32:ASN:ND2	30:D:301:HOH:O	2.29	0.44
1:N:87:ILE:O	1:N:173:PRO:HD3	2.17	0.44
1:N:194:LEU:HD22	1:N:285:PHE:HE2	1.81	0.44
3:P:106:LEU:HA	3:P:106:LEU:HD23	1.71	0.44
3:P:109:THR:O	3:P:112:LEU:HB2	2.17	0.44
26:T:104:CDL:H231	26:T:104:CDL:H511	1.99	0.44
1:N:115:SER:O	1:N:121:GLY:HA2	2.17	0.44
1:N:417:MET:CE	30:N:841:HOH:O	2.65	0.44
3:P:146:TRP:CZ2	7:T:17:ARG:HG3	2.52	0.44
19:O:302:TGL:H162	19:O:302:TGL:HC81	1.68	0.44
1:A:75:ILE:O	1:A:79:GLY:HA3	2.17	0.44
1:A:309:THR:HG22	14:A:602:HEA:HMB2	1.99	0.44
6:F:39:THR:HG23	21:F:102:EDO:H12	1.98	0.44
12:Y:4:GLU:HB3	12:Y:9:LYS:HB3	1.99	0.44
2:B:1:FME:HE1	2:B:133:LEU:HD13	1.99	0.44
3:C:231:HIS:HB2	30:F:211:HOH:O	2.17	0.44
6:F:62:CYS:HB3	6:F:85:CYS:HB3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:G:2:SER:HA	27:G:103:PEK:H331	2.00	0.44
26:G:102:CDL:H852	27:G:103:PEK:H372	1.99	0.44
1:N:489:THR:HA	6:S:71:TRP:O	2.17	0.44
19:O:302:TGL:HB92	19:O:302:TGL:H281	2.00	0.44
3:C:129:VAL:O	3:C:130:PRO:C	2.59	0.44
7:G:3:ALA:O	7:G:4:ALA:CB	2.66	0.44
14:N:601:HEA:HHC	14:N:601:HEA:H11	1.85	0.44
2:O:102:HIS:CE1	2:O:157:GLU:OE2	2.71	0.44
6:S:70:ILE:HG13	6:S:84:SER:HB3	1.98	0.44
1:A:177:SER:HB3	1:A:180:GLN:HE22	1.83	0.44
10:W:12:PHE:O	10:W:23:LYS:HE2	2.16	0.44
8:H:46:LYS:O	8:H:46:LYS:HD3	2.18	0.44
2:O:104:TRP:CG	2:O:203:ASN:HB2	2.53	0.44
1:N:513:LEU:HD23	1:N:513:LEU:HA	1.80	0.44
1:N:12:HIS:CE1	1:N:13:LYS:HG2	2.53	0.43
2:O:59:GLN:O	2:O:60:GLU:HG3	2.17	0.43
3:P:103:HIS:ND1	23:P:306:CHD:O26	2.51	0.43
19:Q:201:TGL:HB51	19:Q:201:TGL:HA21	2.00	0.43
6:F:55:LYS:HA	6:F:74:LEU:O	2.17	0.43
6:F:95:GLN:HG3	6:F:96:LEU:H	1.82	0.43
6:F:95:GLN:CG	6:F:96:LEU:N	2.80	0.43
8:U:39:CYS:O	8:U:43:MET:HG2	2.18	0.43
1:A:312:ILE:O	1:A:312:ILE:HG22	2.17	0.43
4:Q:77:GLU:OE2	19:Q:201:TGL:HB21	2.18	0.43
1:A:203:ALA:HB1	27:G:101:PEK:H383	2.00	0.43
12:L:2:HIS:CG	12:L:3:TYR:H	2.36	0.43
3:P:55:TYR:HA	26:P:304:CDL:H552	2.00	0.43
1:A:344:PHE:CD1	1:A:344:PHE:C	2.97	0.43
1:A:488:THR:HB	1:A:495:LEU:HD13	1.99	0.43
7:G:25:LEU:HD23	7:G:25:LEU:HA	1.83	0.43
12:L:32:GLY:N	30:L:201:HOH:O	2.51	0.43
1:N:466:MET:HG2	13:Z:26:PHE:CG	2.54	0.43
10:W:30:ILE:O	10:W:34:VAL:HG23	2.18	0.43
1:N:309:THR:CG2	14:N:602:HEA:HMB2	2.48	0.43
3:P:22:LEU:HD23	3:P:22:LEU:HA	1.83	0.43
3:P:67:PHE:HE2	26:P:304:CDL:H1	1.83	0.43
9:V:55:ASP:HA	30:V:110:HOH:O	2.18	0.43
1:A:302:ARG:NH2	8:H:23:GLN:HE21	2.17	0.43
1:A:381[B]:LEU:HD13	14:A:602:HEA:HBC2	2.01	0.43
2:B:86:MET:C	2:B:88:ASP:H	2.26	0.43
6:F:55:LYS:NZ	30:F:202:HOH:O	2.37	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:240:HIS:CD2	1:N:240:HIS:C	2.97	0.43
3:P:259:TRP:HB2	30:P:481:HOH:O	2.17	0.43
10:W:33:ARG:CZ	23:W:101:CHD:H212	2.49	0.43
10:W:33:ARG:HD3	30:W:208:HOH:O	2.18	0.43
1:A:467:LEU:O	1:A:471:ILE:HG13	2.19	0.43
2:B:146:MET:SD	2:B:189:PRO:HB3	2.59	0.43
1:N:501:PRO:HD2	1:N:504:THR:CG2	2.49	0.43
2:O:92:ASN:HA	2:O:93:PRO:HD2	1.96	0.43
3:P:30:GLY:HA2	3:P:42:LEU:HB3	2.00	0.43
23:P:305:CHD:H232	23:P:305:CHD:C16	2.46	0.43
7:T:3:ALA:O	7:T:4:ALA:HB2	2.19	0.43
2:B:196:CYS:HB2	2:B:207:MET:HG3	2.00	0.43
1:A:440:TYR:OH	2:B:195:GLN:HB3	2.19	0.42
2:B:86:MET:C	2:B:88:ASP:N	2.77	0.42
3:C:110:PRO:HB3	8:H:30:TRP:CE3	2.54	0.42
30:E:239:HOH:O	9:I:12:LEU:HD12	2.19	0.42
2:O:116:LEU:HD11	2:O:226:MET:HG2	2.01	0.42
3:P:129:VAL:HG11	3:P:180:GLU:CG	2.49	0.42
4:Q:86:MET:O	11:X:25:CYS:HB2	2.19	0.42
1:A:106:PRO:N	1:A:107:PRO:HD2	2.33	0.42
14:A:602:HEA:H243	2:B:69:PRO:HB3	2.01	0.42
20:A:609:PGV:H02	20:A:609:PGV:O14	2.18	0.42
19:Y:101:TGL:H222	19:Y:101:TGL:HA92	1.47	0.42
3:C:223:LEU:HD21	23:C:304:CHD:H183	2.01	0.42
27:C:306:PEK:H032	7:G:17:ARG:NH2	2.34	0.42
4:D:138:TRP:CH2	11:K:50:PRO:HG2	2.55	0.42
5:E:9:GLU:H	5:E:9:GLU:CD	2.27	0.42
1:N:390:MET:HA	1:N:390:MET:CE	2.40	0.42
4:Q:7:LYS:O	4:Q:8:SER:HB2	2.17	0.42
10:W:2:GLU:OE1	10:W:4:ARG:NH2	2.52	0.42
14:A:601:HEA:H271	14:A:601:HEA:H212	1.80	0.42
2:B:78:LEU:HD12	2:B:78:LEU:HA	1.64	0.42
26:G:102:CDL:H392	2:O:78:LEU:HD12	2.00	0.42
4:Q:78:TRP:N	19:Q:201:TGL:HB22	2.33	0.42
1:A:5:ARG:HA	21:A:612:EDO:O1	2.19	0.42
1:A:344:PHE:CD2	1:A:384[A]:GLY:O	2.73	0.42
3:C:187:THR:HG22	27:G:101:PEK:H051	2.00	0.42
6:F:37:LYS:HA	6:F:37:LYS:HD2	1.70	0.42
8:H:46:LYS:C	8:H:48:GLY:N	2.77	0.42
1:N:482:VAL:HG13	13:Z:1:ILE:HD11	2.01	0.42
2:O:50:LEU:HD23	2:O:50:LEU:HA	1.89	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
10:W:50:LEU:O	10:W:50:LEU:HD22	2.20	0.42
1:N:366:VAL:N	30:N:705:HOH:O	2.52	0.42
3:P:157:LYS:NZ	27:T:103:PEK:H051	2.34	0.42
7:T:11:TPO:O	7:T:11:TPO:CG2	2.68	0.42
1:A:399:LEU:HB2	1:A:494:TRP:CZ3	2.54	0.42
1:A:495:LEU:HD12	1:A:495:LEU:HA	1.82	0.42
1:N:459:PHE:O	1:N:462:LEU:HB3	2.20	0.42
1:A:431:LEU:HD21	1:A:450:TRP:HB2	2.01	0.42
26:G:102:CDL:H431	2:O:77:ALA:HB3	2.00	0.42
26:T:104:CDL:H231	26:T:104:CDL:C51	2.49	0.42
12:Y:22:LEU:O	12:Y:26:THR:HB	2.20	0.42
2:B:145:PRO:HA	2:B:214:VAL:O	2.19	0.42
6:F:49:VAL:HA	6:F:50:PRO:HD3	1.89	0.42
1:N:43:GLN:HG2	4:Q:104:TYR:CE1	2.55	0.42
14:N:602:HEA:HHA	14:N:602:HEA:HAD2	1.80	0.42
4:Q:121:LYS:HG2	11:X:52:GLU:HA	2.02	0.42
3:C:193:TYR:CD1	3:C:193:TYR:C	2.98	0.42
8:H:46:LYS:C	8:H:48:GLY:H	2.27	0.42
2:O:164:ALA:O	2:O:194:GLY:HA3	2.20	0.42
3:P:33:MET:HG2	30:P:473:HOH:O	2.20	0.42
8:U:58:ARG:HA	8:U:58:ARG:HD2	1.82	0.42
3:C:125:ASN:HB3	3:C:128:GLU:HG3	2.01	0.41
3:C:223:LEU:HD23	3:C:223:LEU:HA	1.87	0.41
7:T:23:LEU:HB2	21:T:105:EDO:H12	2.02	0.41
2:B:68:LEU:CB	2:B:69:PRO:HD3	2.50	0.41
1:N:113:LEU:O	1:N:114:ALA:C	2.63	0.41
1:N:339:MET:HE1	1:N:411:LYS:HD3	2.02	0.41
14:N:601:HEA:HHD	30:N:759:HOH:O	2.19	0.41
27:T:102:PEK:H381	26:T:104:CDL:H872	1.99	0.41
1:A:344:PHE:HD2	1:A:384[A]:GLY:O	2.03	0.41
3:C:146:TRP:CZ2	7:G:17:ARG:HG3	2.54	0.41
3:C:149:HIS:HA	3:C:152:MET:HE3	2.01	0.41
7:G:44:ARG:HD2	7:G:74:ARG:O	2.20	0.41
12:L:25:MET:HG2	19:L:101:TGL:HA61	2.02	0.41
3:C:105:SER:HA	3:C:116:TRP:CE3	2.56	0.41
5:E:23:ASP:OD1	5:E:23:ASP:N	2.38	0.41
6:F:52:ILE:HG23	6:F:96:LEU:HD13	2.02	0.41
7:G:9:GLY:O	1:N:177:SER:HA	2.21	0.41
1:A:261:TYR:HB2	1:A:338:MET:HE2	2.02	0.41
2:B:131:GLY:HA2	4:D:119:GLN:HA	2.03	0.41
4:Q:102:TYR:CE1	13:Z:35:TYR:HE2	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
13:Z:36:HIS:HD2	13:Z:39:ASN:ND2	2.18	0.41
1:A:449:MET:HE3	1:A:449:MET:HB2	1.95	0.41
7:T:40:GLY:O	7:T:41:HIS:C	2.64	0.41
19:Y:101:TGL:H321	19:Y:101:TGL:H362	2.01	0.41
2:B:10:GLN:HB3	30:B:487:HOH:O	2.20	0.41
26:G:102:CDL:H521	26:G:102:CDL:H242	2.02	0.41
2:B:129:LYS:HE3	30:B:471:HOH:O	2.21	0.41
5:R:36:LEU:HD21	5:R:47:ILE:HG21	2.02	0.41
12:Y:11:ILE:HD12	12:Y:13:PHE:CE1	2.55	0.41
19:Y:101:TGL:HC22	19:Y:101:TGL:HC52	1.93	0.41
1:A:379:TYR:O	1:A:383[B]:MET:HB2	2.21	0.41
19:A:607:TGL:H222	19:A:607:TGL:H362	2.02	0.41
2:B:37:LEU:HB2	9:I:28:SER:OG	2.21	0.41
24:B:303:PSC:H212	24:B:303:PSC:H02	2.02	0.41
3:C:85:LEU:HD21	27:T:102:PEK:H281	2.03	0.41
19:D:201:TGL:HA62	19:D:201:TGL:HB82	2.03	0.41
5:E:31:LYS:HE2	5:E:35:THR:OG1	2.21	0.41
1:N:66:ILE:HG13	30:N:759:HOH:O	2.21	0.41
19:O:302:TGL:C28	19:O:302:TGL:CB9	2.99	0.41
26:T:104:CDL:H252	26:T:104:CDL:H761	2.01	0.41
23:W:101:CHD:H151	23:W:101:CHD:O7	2.20	0.41
1:A:282:PHE:HZ	26:T:104:CDL:H752	1.86	0.41
2:B:171:LYS:HB2	2:B:171:LYS:HE3	1.92	0.41
2:O:82:ARG:O	2:O:83:ILE:C	2.64	0.41
1:A:344:PHE:HD2	1:A:384[B]:GLY:O	2.04	0.40
2:B:28:LEU:HD12	2:B:28:LEU:HA	1.89	0.40
2:O:121:TYR:O	2:O:138:VAL:HA	2.20	0.40
3:P:146:TRP:HB2	7:T:16:TRP:HB3	2.02	0.40
5:R:21:LYS:HA	5:R:22:PRO:HD3	1.94	0.40
5:R:80:GLU:CD	5:R:80:GLU:N	2.79	0.40
7:T:31:CYS:SG	26:T:104:CDL:H512	2.61	0.40
1:A:22:PHE:CE1	30:A:867:HOH:O	2.07	0.40
3:C:22:LEU:HD23	3:C:22:LEU:HA	1.95	0.40
1:N:62:ALA:HA	30:N:759:HOH:O	2.22	0.40
23:C:304:CHD:H112	23:C:304:CHD:H12A	1.95	0.40
4:D:13:LEU:HD23	4:D:13:LEU:HA	1.93	0.40
5:E:12:ASP:HA	5:E:47:ILE:HD11	2.02	0.40
19:L:101:TGL:HC21	30:L:224:HOH:O	2.21	0.40
1:N:76:GLY:O	1:N:80:ASN:HB2	2.21	0.40
1:N:172:LYS:HB2	1:N:172:LYS:HE2	1.87	0.40
1:N:265:LYS:HD2	30:O:483:HOH:O	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:O:42:ILE:O	2:O:46:LEU:HG	2.21	0.40
2:B:200:CYS:SG	2:B:204:HIS:HA	2.61	0.40
7:G:10:GLY:HA3	1:N:177:SER:HB2	2.03	0.40
4:Q:118:LYS:HG2	11:X:51:LYS:HB3	2.03	0.40
10:W:33:ARG:NH2	23:W:101:CHD:H212	2.37	0.40
1:A:183:LEU:HA	1:A:183:LEU:HD23	1.83	0.40
2:B:1:FME:HB2	2:B:1:FME:HCN	1.76	0.40
24:B:303:PSC:H241	24:B:303:PSC:H21	2.03	0.40
4:D:48:TRP:HA	4:D:51:LEU:HD22	2.04	0.40
23:P:306:CHD:H12	23:P:306:CHD:C21	2.47	0.40
7:T:25:LEU:HD23	7:T:25:LEU:HA	1.88	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	519/514 (101%)	503 (97%)	16 (3%)	0	100	100
1	N	512/514 (100%)	493 (96%)	19 (4%)	0	100	100
2	B	225/227 (99%)	215 (96%)	10 (4%)	0	100	100
2	O	225/227 (99%)	217 (96%)	6 (3%)	2 (1%)	14	14
3	C	257/261 (98%)	251 (98%)	5 (2%)	1 (0%)	30	34
3	P	257/261 (98%)	253 (98%)	4 (2%)	0	100	100
4	D	142/147 (97%)	139 (98%)	3 (2%)	0	100	100
4	Q	142/147 (97%)	137 (96%)	4 (3%)	1 (1%)	18	19
5	E	103/109 (94%)	101 (98%)	2 (2%)	0	100	100
5	R	103/109 (94%)	101 (98%)	2 (2%)	0	100	100
6	F	96/98 (98%)	89 (93%)	4 (4%)	3 (3%)	3	1

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
6	S	96/98 (98%)	88 (92%)	8 (8%)	0	100	100
7	G	81/85 (95%)	69 (85%)	9 (11%)	3 (4%)	2	1
7	T	81/85 (95%)	69 (85%)	5 (6%)	7 (9%)	0	0
8	H	77/85 (91%)	71 (92%)	4 (5%)	2 (3%)	4	2
8	U	77/85 (91%)	70 (91%)	4 (5%)	3 (4%)	2	1
9	I	72/73 (99%)	71 (99%)	1 (1%)	0	100	100
9	V	73/73 (100%)	71 (97%)	1 (1%)	1 (1%)	9	7
10	J	56/59 (95%)	54 (96%)	2 (4%)	0	100	100
10	W	56/59 (95%)	55 (98%)	1 (2%)	0	100	100
11	K	47/56 (84%)	47 (100%)	0	0	100	100
11	X	47/56 (84%)	42 (89%)	5 (11%)	0	100	100
12	L	44/47 (94%)	44 (100%)	0	0	100	100
12	Y	44/47 (94%)	42 (96%)	2 (4%)	0	100	100
13	M	41/46 (89%)	41 (100%)	0	0	100	100
13	Z	41/46 (89%)	41 (100%)	0	0	100	100
All	All	3514/3614 (97%)	3374 (96%)	117 (3%)	23 (1%)	18	19

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	F	94	HIS
6	F	96	LEU
7	G	4	ALA
7	G	39	SER
7	T	4	ALA
7	T	7	ASP
7	T	8	HIS
7	T	39	SER
9	V	2	THR
8	H	47	GLY
2	O	60	GLU
4	Q	8	SER
7	T	40	GLY
7	T	41	HIS
8	U	48	GLY
7	G	43	GLU
2	O	92	ASN

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Mol	Chain	Res	Type
8	U	51	SER
3	C	38	ASN
6	F	95	GLN
8	H	8	ILE
8	U	8	ILE
7	T	6	GLY

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	432/426 (101%)	423 (98%)	9 (2%)	47	63
1	N	426/426 (100%)	411 (96%)	15 (4%)	32	43
2	B	210/210 (100%)	200 (95%)	10 (5%)	23	30
2	O	210/210 (100%)	194 (92%)	16 (8%)	12	14
3	C	224/226 (99%)	219 (98%)	5 (2%)	45	61
3	P	224/226 (99%)	220 (98%)	4 (2%)	51	68
4	D	128/129 (99%)	121 (94%)	7 (6%)	19	24
4	Q	128/129 (99%)	122 (95%)	6 (5%)	23	31
5	E	92/95 (97%)	90 (98%)	2 (2%)	45	61
5	R	92/95 (97%)	86 (94%)	6 (6%)	15	18
6	F	81/81 (100%)	72 (89%)	9 (11%)	6	6
6	S	81/81 (100%)	72 (89%)	9 (11%)	6	6
7	G	67/68 (98%)	59 (88%)	8 (12%)	5	4
7	T	67/68 (98%)	61 (91%)	6 (9%)	9	9
8	H	71/75 (95%)	60 (84%)	11 (16%)	2	2
8	U	71/75 (95%)	65 (92%)	6 (8%)	10	11
9	I	58/57 (102%)	50 (86%)	8 (14%)	3	3
9	V	59/57 (104%)	52 (88%)	7 (12%)	5	4
10	J	49/50 (98%)	45 (92%)	4 (8%)	10	12

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
10	W	49/50 (98%)	46 (94%)	3 (6%)	17	20
11	K	39/46 (85%)	36 (92%)	3 (8%)	12	13
11	X	39/46 (85%)	37 (95%)	2 (5%)	21	27
12	L	39/40 (98%)	36 (92%)	3 (8%)	12	13
12	Y	39/40 (98%)	37 (95%)	2 (5%)	21	27
13	M	37/38 (97%)	34 (92%)	3 (8%)	11	12
13	Z	37/38 (97%)	33 (89%)	4 (11%)	6	6
All	All	3049/3082 (99%)	2881 (94%)	168 (6%)	19	24

All (168) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	ARG
1	A	169	ILE
1	A	177	SER
1	A	180	GLN
1	A	312	ILE
1	A	369	ASP
1	A	394	VAL
1	A	513	LEU
1	A	514	LYS
2	B	15	PRO
2	B	55	THR
2	B	57	ASP
2	B	61	VAL
2	B	65	TRP
2	B	66	THR
2	B	68	LEU
2	B	75	LEU
2	B	78	LEU
2	B	217	LYS
3	C	23	SER
3	C	51	MET
3	C	127	LEU
3	C	159	MET
3	C	246	ASP
4	D	4	SER
4	D	8	SER
4	D	45	LYS
4	D	51	LEU

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Mol	Chain	Res	Type
4	D	58	GLU
4	D	121	LYS
4	D	147	LYS
5	E	41	LEU
5	E	70	VAL
6	F	37	LYS
6	F	43	LYS
6	F	48	LEU
6	F	53	THR
6	F	80	GLN
6	F	87	THR
6	F	94	HIS
6	F	95	GLN
6	F	96	LEU
7	G	2	SER
7	G	17	ARG
7	G	33	LEU
7	G	35	SER
7	G	36	TRP
7	G	37	LEU
7	G	42	ARG
7	G	74	ARG
8	H	7	LYS
8	H	9	LYS
8	H	23	GLN
8	H	29	CYS
8	H	46	LYS
8	H	50	VAL
8	H	52	VAL
8	H	60	TYR
8	H	61	LYS
8	H	70	SER
8	H	84	LYS
9	I	2	THR
9	I	8	GLN
9	I	18	ARG
9	I	25	PHE
9	I	28	SER
9	I	29	LEU
9	I	53	ASN
9	I	68	ILE
10	J	29	ASN

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Mol	Chain	Res	Type
10	J	31	LEU
10	J	50	LEU
10	J	58	LYS
11	K	20	SER
11	K	38	ILE
11	K	54	ARG
12	L	5	GLU
12	L	26	THR
12	L	47	LYS
13	M	34	LEU
13	M	42	LYS
13	M	43	SER
1	N	57	VAL
1	N	118	VAL
1	N	158	ILE
1	N	178	GLN
1	N	180	GLN
1	N	265	LYS
1	N	312	ILE
1	N	361	SER
1	N	363	LEU
1	N	369	ASP
1	N	380	VAL
1	N	394	VAL
1	N	486	ASP
1	N	495	LEU
1	N	510	TYR
2	O	33	LEU
2	O	42	ILE
2	O	55	THR
2	O	60	GLU
2	O	68	LEU
2	O	75	LEU
2	O	78	LEU
2	O	88	ASP
2	O	89	GLU
2	O	91	ASN
2	O	94	SER
2	O	157	GLU
2	O	183	THR
2	O	185	MET
2	O	221	LYS

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Mol	Chain	Res	Type
2	O	227	LEU
3	P	3	HIS
3	P	40	MET
3	P	127	LEU
3	P	159	MET
4	Q	6	VAL
4	Q	8	SER
4	Q	65	LYS
4	Q	73	ARG
4	Q	74	SER
4	Q	108	PRO
5	R	6	GLU
5	R	21	LYS
5	R	72	LYS
5	R	80	GLU
5	R	108	LYS
5	R	109	VAL
6	S	26	LYS
6	S	43	LYS
6	S	48	LEU
6	S	53	THR
6	S	54	ASN
6	S	63	GLU
6	S	80	GLN
6	S	87	THR
6	S	98	HIS
7	T	17	ARG
7	T	33	LEU
7	T	35	SER
7	T	36	TRP
7	T	37	LEU
7	T	42	ARG
8	U	7	LYS
8	U	8	ILE
8	U	10	ASN
8	U	12	GLN
8	U	46	LYS
8	U	60	TYR
9	V	8	GLN
9	V	21	ILE
9	V	29	LEU
9	V	37[A]	PHE

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Mol	Chain	Res	Type
9	V	37[B]	PHE
9	V	58	LYS
9	V	65	LYS
10	W	4	ARG
10	W	27	THR
10	W	50	LEU
11	X	47	ARG
11	X	54	ARG
12	Y	11	ILE
12	Y	26	THR
13	Z	19	LEU
13	Z	34	LEU
13	Z	37	LEU
13	Z	43	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (63) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	178	GLN
1	A	180	GLN
1	A	512	ASN
2	B	10	GLN
2	B	24	HIS
2	B	181	GLN
2	B	203	ASN
3	C	3	HIS
3	C	50	ASN
3	C	68	GLN
3	C	161	GLN
4	D	29	HIS
4	D	37	GLN
4	D	101	HIS
4	D	109	HIS
5	E	94	ASN
6	F	54	ASN
6	F	80	GLN
6	F	88	HIS
6	F	95	GLN
7	G	34	ASN
7	G	71	HIS
7	G	76	ASN
8	H	23	GLN

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Mol	Chain	Res	Type
8	H	31	GLN
8	H	32	ASN
9	I	8	GLN
9	I	53	ASN
10	J	29	ASN
10	J	57	HIS
11	K	15	ASN
11	K	35	GLN
13	M	39	ASN
1	N	4	ASN
1	N	178	GLN
1	N	180	GLN
1	N	406	ASN
1	N	422	ASN
1	N	512	ASN
2	O	10	GLN
2	O	91	ASN
2	O	181	GLN
2	O	203	ASN
3	P	68	GLN
3	P	161	GLN
4	Q	29	HIS
4	Q	37	GLN
4	Q	109	HIS
4	Q	119	GLN
5	R	94	ASN
6	S	54	ASN
6	S	80	GLN
7	T	8	HIS
7	T	38	HIS
7	T	76	ASN
8	U	10	ASN
8	U	23	GLN
8	U	28	ASN
8	U	31	GLN
8	U	32	ASN
9	V	8	GLN
11	X	35	GLN
13	Z	39	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

8 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
9	SAC	I	1	9	7,8,9	1.45	1 (14%)	7,9,11	1.10	0
2	FME	O	1	2	8,9,10	0.97	0	8,9,11	3.94	2 (25%)
1	FME	A	1	1	8,9,10	0.87	0	8,9,11	5.34	4 (50%)
9	SAC	V	1	9	7,8,9	1.64	1 (14%)	7,9,11	1.61	1 (14%)
7	TPO	T	11	7	8,10,11	1.59	2 (25%)	10,14,16	1.19	1 (10%)
1	FME	N	1	1	8,9,10	0.47	0	8,9,11	3.11	3 (37%)
7	TPO	G	11	7	8,10,11	1.79	2 (25%)	10,14,16	1.01	0
2	FME	B	1	2	8,9,10	0.99	0	8,9,11	5.48	2 (25%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	SAC	I	1	9	-	4/7/8/10	-
2	FME	O	1	2	-	2/7/9/11	-
1	FME	A	1	1	-	4/7/9/11	-
9	SAC	V	1	9	-	6/7/8/10	-
7	TPO	T	11	7	-	3/9/11/13	-
1	FME	N	1	1	-	4/7/9/11	-
7	TPO	G	11	7	-	3/9/11/13	-
2	FME	B	1	2	-	3/7/9/11	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	V	1	SAC	CA-N	4.06	1.52	1.46
9	I	1	SAC	CA-N	3.55	1.51	1.46
7	G	11	TPO	P-O1P	2.98	1.59	1.50
7	T	11	TPO	P-O1P	2.91	1.59	1.50
7	G	11	TPO	P-OG1	2.59	1.64	1.59
7	T	11	TPO	P-OG1	2.10	1.63	1.59

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1	FME	CA-N-CN	-14.94	99.85	122.82
1	A	1	FME	CA-N-CN	-13.87	101.50	122.82
2	O	1	FME	CA-N-CN	-10.10	107.28	122.82
1	N	1	FME	CA-N-CN	-7.45	111.36	122.82
2	O	1	FME	C-CA-N	3.71	116.67	109.50
1	A	1	FME	O1-CN-N	3.46	134.25	125.32
1	N	1	FME	CB-CA-N	3.43	116.76	110.52
9	V	1	SAC	C-CA-N	3.41	116.09	109.50
2	B	1	FME	C-CA-N	3.17	115.62	109.50
1	A	1	FME	CG-CB-CA	-2.73	104.52	112.87
7	T	11	TPO	O3P-P-OG1	2.54	115.74	105.85
1	A	1	FME	CB-CA-N	2.47	115.01	110.52
1	N	1	FME	O-C-CA	-2.37	118.67	124.77

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1	FME	O1-CN-N-CA
1	A	1	FME	N-CA-CB-CG
2	B	1	FME	O1-CN-N-CA
7	G	11	TPO	N-CA-CB-OG1
9	I	1	SAC	C2A-C1A-N-CA
9	I	1	SAC	OAC-C1A-N-CA
1	N	1	FME	O1-CN-N-CA
1	N	1	FME	N-CA-CB-CG
2	O	1	FME	O1-CN-N-CA
7	T	11	TPO	N-CA-CB-CG2
7	T	11	TPO	N-CA-CB-OG1
7	T	11	TPO	C-CA-CB-CG2
9	V	1	SAC	C-CA-N-C1A

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Mol	Chain	Res	Type	Atoms
9	V	1	SAC	CB-CA-N-C1A
9	V	1	SAC	C2A-C1A-N-CA
9	V	1	SAC	OAC-C1A-N-CA
9	V	1	SAC	N-CA-CB-OG
2	B	1	FME	CB-CG-SD-CE
7	G	11	TPO	CG2-CB-OG1-P
1	A	1	FME	CB-CG-SD-CE
9	V	1	SAC	C-CA-CB-OG
9	I	1	SAC	C-CA-N-C1A
1	A	1	FME	C-CA-CB-CG
1	N	1	FME	C-CA-CB-CG
2	B	1	FME	CB-CA-N-CN
2	O	1	FME	CB-CA-N-CN
9	I	1	SAC	CB-CA-N-C1A
1	N	1	FME	CB-CG-SD-CE
7	G	11	TPO	CB-OG1-P-O3P

There are no ring outliers.

4 monomers are involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	O	1	FME	1	0
7	T	11	TPO	1	0
1	N	1	FME	1	0
2	B	1	FME	4	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 77 ligands modelled in this entry, 8 are monoatomic and 2 are unknown - leaving 67 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
14	HEA	N	602	1	67,67,67	1.93	19 (28%)	81,103,103	1.94	25 (30%)
18	CMO	A	606	15	0,1,1	-	-	-		
18	CMO	N	606	-	0,1,1	-	-	-		
22	CUA	O	301	30,2	0,1,1	-	-	-		
23	CHD	J	101	-	32,32,32	0.80	0	51,51,51	1.86	9 (17%)
27	PEK	T	101	-	52,52,52	0.91	3 (5%)	55,57,57	1.01	3 (5%)
14	HEA	A	601	1	67,67,67	2.00	20 (29%)	81,103,103	1.60	16 (19%)
29	DMU	M	101	-	34,34,34	0.46	0	45,45,45	0.94	1 (2%)
19	TGL	L	101	-	62,62,62	1.20	3 (4%)	65,65,65	1.45	8 (12%)
21	EDO	K	102	-	3,3,3	0.44	0	2,2,2	0.35	0
21	EDO	L	102	-	3,3,3	0.39	0	2,2,2	0.35	0
20	PGV	N	607	-	50,50,50	1.04	2 (4%)	53,56,56	1.13	5 (9%)
21	EDO	B	305	-	3,3,3	0.50	0	2,2,2	0.86	0
23	CHD	W	101	-	32,32,32	0.90	0	51,51,51	4.03	25 (49%)
27	PEK	T	102	-	52,52,52	1.12	2 (3%)	55,57,57	0.99	5 (9%)
19	TGL	O	302	-	62,62,62	1.12	3 (4%)	65,65,65	1.26	7 (10%)
27	PEK	C	306	-	52,52,52	1.18	3 (5%)	55,57,57	1.18	5 (9%)
21	EDO	C	311	-	3,3,3	0.71	0	2,2,2	0.38	0
20	PGV	C	307	-	50,50,50	1.05	2 (4%)	53,56,56	0.93	3 (5%)
20	PGV	C	302	-	50,50,50	0.89	2 (4%)	53,56,56	0.96	4 (7%)
26	CDL	C	303	-	99,99,99	1.48	12 (12%)	105,111,111	1.25	9 (8%)
19	TGL	Y	101	-	62,62,62	1.15	3 (4%)	65,65,65	1.26	6 (9%)
21	EDO	A	611	-	3,3,3	0.38	0	2,2,2	1.09	0
20	PGV	P	303	-	50,50,50	0.87	4 (8%)	53,56,56	1.19	7 (13%)
21	EDO	N	610	-	3,3,3	0.39	0	2,2,2	0.24	0
14	HEA	A	602	1	67,67,67	1.85	19 (28%)	81,103,103	2.20	27 (33%)
21	EDO	F	103	-	3,3,3	0.50	0	2,2,2	0.43	0
27	PEK	T	103	-	52,52,52	1.13	2 (3%)	55,57,57	0.96	2 (3%)
24	PSC	B	303	-	51,51,51	1.21	3 (5%)	57,59,59	1.14	4 (7%)
21	EDO	T	105	-	3,3,3	0.45	0	2,2,2	0.30	0
21	EDO	P	307	-	3,3,3	0.51	0	2,2,2	0.15	0
20	PGV	N	608	-	50,50,50	0.88	2 (4%)	53,56,56	1.10	4 (7%)
21	EDO	P	308	-	3,3,3	0.56	0	2,2,2	0.19	0
26	CDL	P	304	-	99,99,99	1.50	12 (12%)	105,111,111	1.29	7 (6%)
19	TGL	D	201	-	62,62,62	1.31	3 (4%)	65,65,65	1.09	6 (9%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
21	EDO	F	102	-	3,3,3	0.51	0	2,2,2	0.15	0
21	EDO	O	304	-	3,3,3	0.36	0	2,2,2	0.70	0
26	CDL	G	102	-	99,99,99	1.50	12 (12%)	105,111,111	1.23	9 (8%)
23	CHD	G	104	-	32,32,32	1.18	2 (6%)	51,51,51	1.52	11 (21%)
21	EDO	A	613	-	3,3,3	0.53	0	2,2,2	0.01	0
27	PEK	G	103	-	52,52,52	1.17	2 (3%)	55,57,57	1.09	5 (9%)
19	TGL	Q	201	-	62,62,62	1.15	3 (4%)	65,65,65	1.07	5 (7%)
20	PGV	A	608	-	50,50,50	0.88	2 (4%)	53,56,56	1.48	4 (7%)
21	EDO	B	304	-	3,3,3	0.32	0	2,2,2	0.65	0
26	CDL	T	104	-	99,99,99	1.45	12 (12%)	105,111,111	1.18	5 (4%)
21	EDO	A	610	-	3,3,3	0.46	0	2,2,2	0.30	0
14	HEA	N	601	1	67,67,67	2.06	25 (37%)	81,103,103	1.38	12 (14%)
21	EDO	C	309	-	3,3,3	0.49	0	2,2,2	0.27	0
23	CHD	C	305	-	32,32,32	0.88	1 (3%)	51,51,51	1.66	11 (21%)
23	CHD	P	306	-	32,32,32	0.84	0	51,51,51	1.61	11 (21%)
20	PGV	P	301	-	50,50,50	1.08	2 (4%)	53,56,56	1.12	4 (7%)
29	DMU	Z	101	-	34,34,34	0.49	0	45,45,45	0.86	2 (4%)
19	TGL	A	607	-	62,62,62	1.20	4 (6%)	65,65,65	1.71	8 (12%)
21	EDO	A	612	-	3,3,3	0.43	0	2,2,2	0.35	0
23	CHD	P	305	-	32,32,32	0.62	0	51,51,51	1.74	13 (25%)
23	CHD	C	304	-	32,32,32	0.58	0	51,51,51	1.95	17 (33%)
23	CHD	B	302	-	32,32,32	0.76	0	51,51,51	1.46	11 (21%)
24	PSC	O	303	-	51,51,51	1.20	3 (5%)	57,59,59	1.07	4 (7%)
21	EDO	K	101	-	3,3,3	0.49	0	2,2,2	0.34	0
21	EDO	N	609	-	3,3,3	0.39	0	2,2,2	0.67	0
21	EDO	C	308	-	3,3,3	0.39	0	2,2,2	0.42	0
21	EDO	G	105	-	3,3,3	0.39	0	2,2,2	0.58	0
21	EDO	H	101	-	3,3,3	0.50	0	2,2,2	0.24	0
27	PEK	G	101	-	52,52,52	0.88	3 (5%)	55,57,57	1.38	5 (9%)
20	PGV	A	609	-	50,50,50	1.05	2 (4%)	53,56,56	1.19	6 (11%)
21	EDO	C	310	-	3,3,3	0.54	0	2,2,2	0.39	0
22	CUA	B	301	2	0,1,1	-	-	-	-	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	HEA	N	602	1	-	8/36/76/76	-
23	CHD	J	101	-	-	8/9/74/74	0/4/4/4
29	DMU	M	101	-	-	7/19/59/59	0/2/2/2
27	PEK	T	101	-	-	23/56/56/56	-
14	HEA	A	601	1	-	2/36/76/76	-
19	TGL	L	101	-	-	40/65/65/65	-
21	EDO	K	102	-	-	1/1/1/1	-
21	EDO	L	102	-	-	1/1/1/1	-
20	PGV	N	607	-	-	27/55/55/55	-
21	EDO	B	305	-	-	1/1/1/1	-
23	CHD	W	101	-	-	7/9/74/74	0/4/4/4
27	PEK	T	102	-	-	27/56/56/56	-
19	TGL	O	302	-	-	36/65/65/65	-
27	PEK	C	306	-	-	28/56/56/56	-
21	EDO	C	311	-	-	0/1/1/1	-
20	PGV	C	307	-	-	27/55/55/55	-
20	PGV	C	302	-	-	10/55/55/55	-
26	CDL	C	303	-	-	57/110/110/110	-
19	TGL	Y	101	-	-	41/65/65/65	-
21	EDO	A	611	-	-	1/1/1/1	-
20	PGV	P	303	-	-	20/55/55/55	-
21	EDO	N	610	-	-	0/1/1/1	-
14	HEA	A	602	1	-	2/36/76/76	-
21	EDO	F	103	-	-	0/1/1/1	-
27	PEK	T	103	-	-	31/56/56/56	-
24	PSC	B	303	-	-	37/55/55/55	-
21	EDO	T	105	-	-	1/1/1/1	-
21	EDO	P	307	-	-	0/1/1/1	-
20	PGV	N	608	-	-	14/55/55/55	-
21	EDO	P	308	-	-	1/1/1/1	-
26	CDL	P	304	-	-	64/110/110/110	-
19	TGL	D	201	-	-	37/65/65/65	-
21	EDO	F	102	-	-	0/1/1/1	-
21	EDO	O	304	-	-	1/1/1/1	-
26	CDL	G	102	-	-	64/110/110/110	-
23	CHD	G	104	-	-	3/9/74/74	0/4/4/4
21	EDO	A	613	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
27	PEK	G	103	-	-	26/56/56/56	-
19	TGL	Q	201	-	-	35/65/65/65	-
20	PGV	A	608	-	-	15/55/55/55	-
21	EDO	B	304	-	-	1/1/1/1	-
26	CDL	T	104	-	-	58/110/110/110	-
21	EDO	A	610	-	-	0/1/1/1	-
14	HEA	N	601	1	-	4/36/76/76	-
21	EDO	C	309	-	-	0/1/1/1	-
23	CHD	C	305	-	-	2/9/74/74	0/4/4/4
23	CHD	P	306	-	-	1/9/74/74	0/4/4/4
20	PGV	P	301	-	-	25/55/55/55	-
29	DMU	Z	101	-	-	5/19/59/59	0/2/2/2
19	TGL	A	607	-	-	42/65/65/65	-
21	EDO	A	612	-	-	1/1/1/1	-
23	CHD	P	305	-	-	2/9/74/74	0/4/4/4
23	CHD	C	304	-	-	2/9/74/74	0/4/4/4
23	CHD	B	302	-	-	3/9/74/74	0/4/4/4
24	PSC	O	303	-	-	34/55/55/55	-
21	EDO	K	101	-	-	1/1/1/1	-
21	EDO	N	609	-	-	0/1/1/1	-
21	EDO	C	308	-	-	1/1/1/1	-
21	EDO	G	105	-	-	0/1/1/1	-
21	EDO	H	101	-	-	0/1/1/1	-
27	PEK	G	101	-	-	18/56/56/56	-
20	PGV	A	609	-	-	37/55/55/55	-
21	EDO	C	310	-	-	0/1/1/1	-

All (192) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
19	L	101	TGL	OG2-CB1	5.61	1.50	1.34
19	D	201	TGL	OG2-CB1	5.59	1.50	1.34
26	P	304	CDL	OA6-CA5	5.57	1.50	1.34
14	N	602	HEA	FE-ND	5.54	2.12	1.94
19	A	607	TGL	OG2-CB1	5.51	1.49	1.34
26	G	102	CDL	OB8-CB7	5.51	1.49	1.33
26	G	102	CDL	OB6-CB5	5.45	1.49	1.34
27	C	306	PEK	O01-C1	5.33	1.49	1.34
19	Y	101	TGL	OG2-CB1	5.25	1.49	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
27	G	103	PEK	O03-C21	5.23	1.48	1.33
19	D	201	TGL	OG1-CA1	5.16	1.48	1.33
27	C	306	PEK	O03-C21	5.15	1.48	1.33
27	T	103	PEK	O01-C1	5.15	1.48	1.34
20	C	307	PGV	O01-C1	5.11	1.48	1.34
24	B	303	PSC	O03-C19	5.08	1.48	1.33
26	T	104	CDL	OB6-CB5	5.04	1.48	1.34
19	D	201	TGL	OG3-CC1	4.99	1.47	1.33
27	T	102	PEK	O03-C21	4.98	1.47	1.33
19	Q	201	TGL	OG1-CA1	4.98	1.47	1.33
26	C	303	CDL	OB8-CB7	4.97	1.47	1.33
14	N	601	HEA	FE-ND	4.96	2.10	1.94
19	A	607	TGL	OG1-CA1	4.96	1.47	1.33
26	T	104	CDL	OB8-CB7	4.90	1.47	1.33
20	P	301	PGV	O03-C19	4.89	1.47	1.33
27	T	103	PEK	O03-C21	4.87	1.47	1.33
26	P	304	CDL	OA8-CA7	4.87	1.47	1.33
26	C	303	CDL	OB6-CB5	4.85	1.48	1.34
19	O	302	TGL	OG2-CB1	4.85	1.48	1.34
19	L	101	TGL	OG3-CC1	4.83	1.47	1.33
24	O	303	PSC	O01-C1	4.83	1.47	1.34
14	A	601	HEA	CHB-C4A	4.81	1.47	1.38
24	O	303	PSC	O03-C19	4.78	1.47	1.33
14	N	601	HEA	FE-NB	4.77	2.09	1.94
26	C	303	CDL	OA8-CA7	4.75	1.47	1.33
26	T	104	CDL	OA8-CA7	4.75	1.47	1.33
19	O	302	TGL	OG1-CA1	4.74	1.47	1.33
26	C	303	CDL	OA6-CA5	4.72	1.47	1.34
27	G	103	PEK	O01-C1	4.72	1.47	1.34
19	Y	101	TGL	OG3-CC1	4.69	1.47	1.33
26	P	304	CDL	OB6-CB5	4.68	1.47	1.34
14	A	601	HEA	FE-NB	4.68	2.09	1.94
26	P	304	CDL	OB8-CB7	4.68	1.47	1.33
20	N	607	PGV	O03-C19	4.67	1.47	1.33
19	L	101	TGL	OG1-CA1	4.67	1.47	1.33
14	A	602	HEA	FE-ND	4.65	2.09	1.94
14	A	601	HEA	FE-ND	4.63	2.09	1.94
26	G	102	CDL	OA6-CA5	4.56	1.47	1.34
20	P	301	PGV	O01-C1	4.56	1.47	1.34
19	Q	201	TGL	OG2-CB1	4.55	1.47	1.34
14	N	602	HEA	C4D-C3D	-4.54	1.37	1.45
26	G	102	CDL	OA8-CA7	4.50	1.46	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
20	A	609	PGV	O03-C19	4.48	1.46	1.33
14	N	602	HEA	C1D-ND	-4.46	1.32	1.40
20	A	609	PGV	O01-C1	4.45	1.46	1.34
27	T	102	PEK	O01-C1	4.45	1.46	1.34
24	B	303	PSC	O01-C1	4.45	1.46	1.34
19	O	302	TGL	OG3-CC1	4.40	1.46	1.33
19	Y	101	TGL	OG1-CA1	4.38	1.46	1.33
19	Q	201	TGL	OG3-CC1	4.36	1.46	1.33
14	N	601	HEA	CHA-C1A	4.29	1.46	1.38
20	N	607	PGV	O01-C1	4.23	1.46	1.34
26	T	104	CDL	OA6-CA5	4.17	1.46	1.34
20	C	302	PGV	O03-C19	4.07	1.45	1.33
20	C	307	PGV	O03-C19	4.03	1.45	1.33
14	N	601	HEA	C3A-C4A	-4.02	1.39	1.46
14	A	601	HEA	CHA-C1A	3.86	1.45	1.38
14	A	602	HEA	FE-NB	3.75	2.06	1.94
26	T	104	CDL	C79-C78	-3.74	1.33	1.51
14	A	601	HEA	C4D-C3D	-3.74	1.38	1.45
14	N	602	HEA	FE-NB	3.73	2.06	1.94
24	O	303	PSC	C13-C12	3.71	1.52	1.31
14	A	601	HEA	C1D-ND	-3.69	1.33	1.40
24	B	303	PSC	C13-C12	3.64	1.52	1.31
26	C	303	CDL	C59-C58	-3.64	1.33	1.51
14	N	602	HEA	CHB-C4A	3.63	1.45	1.38
27	T	101	PEK	O03-C21	3.61	1.43	1.33
14	A	602	HEA	C4B-NB	-3.60	1.34	1.40
26	C	303	CDL	C62-C61	-3.59	1.33	1.51
26	P	304	CDL	C22-C21	-3.59	1.33	1.51
20	P	303	PGV	O01-C1	3.58	1.44	1.34
19	A	607	TGL	OG3-CC1	3.58	1.43	1.33
26	P	304	CDL	C79-C78	-3.58	1.34	1.51
26	T	104	CDL	C19-C18	-3.56	1.34	1.51
26	G	102	CDL	C39-C38	-3.56	1.34	1.51
20	A	608	PGV	O03-C19	3.56	1.43	1.33
20	N	608	PGV	O01-C1	3.52	1.44	1.34
26	G	102	CDL	C19-C18	-3.52	1.34	1.51
26	P	304	CDL	C59-C58	-3.51	1.34	1.51
26	G	102	CDL	C79-C78	-3.50	1.34	1.51
26	P	304	CDL	C39-C38	-3.50	1.34	1.51
26	P	304	CDL	C19-C18	-3.49	1.34	1.51
26	T	104	CDL	C59-C58	-3.47	1.34	1.51
26	T	104	CDL	C22-C21	-3.46	1.34	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
20	N	608	PGV	O03-C19	3.46	1.43	1.33
26	C	303	CDL	C79-C78	-3.46	1.34	1.51
26	P	304	CDL	C62-C61	-3.46	1.34	1.51
26	G	102	CDL	C59-C58	-3.45	1.34	1.51
26	T	104	CDL	C39-C38	-3.45	1.34	1.51
14	N	601	HEA	FE-NC	3.44	2.06	1.95
26	C	303	CDL	C82-C81	-3.44	1.34	1.51
26	G	102	CDL	C82-C81	-3.42	1.34	1.51
14	N	602	HEA	CHD-C4C	3.41	1.47	1.39
26	C	303	CDL	C19-C18	-3.40	1.34	1.51
26	T	104	CDL	C62-C61	-3.40	1.34	1.51
26	P	304	CDL	C82-C81	-3.40	1.34	1.51
14	N	601	HEA	C1D-ND	-3.40	1.34	1.40
26	C	303	CDL	C22-C21	-3.40	1.34	1.51
26	C	303	CDL	C39-C38	-3.39	1.34	1.51
27	T	101	PEK	O01-C1	3.39	1.43	1.34
26	P	304	CDL	C42-C41	-3.38	1.35	1.51
14	N	601	HEA	FE-NA	3.38	2.06	1.95
14	N	602	HEA	C1A-C2A	-3.38	1.39	1.45
26	G	102	CDL	C42-C41	-3.37	1.35	1.51
27	G	101	PEK	O03-C21	3.37	1.43	1.33
26	G	102	CDL	C62-C61	-3.37	1.35	1.51
14	N	601	HEA	CHB-C4A	3.36	1.45	1.38
14	A	602	HEA	CHB-C4A	3.36	1.45	1.38
26	T	104	CDL	C42-C41	-3.36	1.35	1.51
26	C	303	CDL	C42-C41	-3.36	1.35	1.51
26	T	104	CDL	C82-C81	-3.34	1.35	1.51
26	G	102	CDL	C22-C21	-3.33	1.35	1.51
14	N	601	HEA	C4B-C3B	-3.32	1.38	1.44
14	A	601	HEA	C4B-NB	-3.31	1.34	1.40
14	A	602	HEA	C1A-C2A	-3.30	1.39	1.45
14	A	601	HEA	FE-NA	3.30	2.06	1.95
27	G	101	PEK	O01-C1	3.28	1.43	1.34
14	A	601	HEA	CHC-C1C	3.27	1.46	1.39
14	A	601	HEA	CHC-C4B	3.20	1.44	1.38
14	A	601	HEA	C3A-C4A	-3.12	1.40	1.46
14	N	601	HEA	CHB-C1B	3.06	1.46	1.39
20	P	303	PGV	O03-C19	3.06	1.42	1.33
14	A	601	HEA	CHD-C1D	3.04	1.44	1.38
20	A	608	PGV	O01-C1	3.01	1.42	1.34
14	A	602	HEA	CHD-C1D	2.92	1.44	1.38
14	A	602	HEA	C1B-NB	-2.92	1.33	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	N	601	HEA	CHD-C1D	2.91	1.44	1.38
14	N	601	HEA	CHA-C4D	2.90	1.45	1.39
14	N	601	HEA	C4B-NB	-2.90	1.35	1.40
14	A	602	HEA	C3A-C4A	-2.88	1.41	1.46
14	A	602	HEA	C1D-ND	-2.86	1.35	1.40
14	A	602	HEA	CMA-C3A	2.82	1.51	1.45
14	N	602	HEA	CHC-C1C	2.78	1.45	1.39
14	N	602	HEA	CHA-C1A	2.77	1.43	1.38
14	N	601	HEA	C1C-NC	-2.77	1.34	1.39
27	T	101	PEK	O01-C02	-2.76	1.40	1.46
14	A	602	HEA	CHC-C1C	2.73	1.45	1.39
27	G	101	PEK	O03-C01	-2.71	1.39	1.45
14	A	602	HEA	CHB-C1B	2.66	1.45	1.39
14	N	602	HEA	C18-C19	2.66	1.39	1.33
14	N	601	HEA	C4D-ND	-2.63	1.33	1.38
14	A	601	HEA	CHD-C4C	2.61	1.45	1.39
14	N	602	HEA	C12-C11	2.61	1.57	1.53
14	N	601	HEA	CHC-C1C	2.61	1.45	1.39
14	A	602	HEA	C4D-ND	-2.60	1.33	1.38
14	A	602	HEA	CHA-C1A	2.60	1.43	1.38
14	N	601	HEA	C4D-C3D	-2.58	1.40	1.45
14	N	602	HEA	CMA-C3A	2.54	1.51	1.45
14	N	602	HEA	FE-NC	2.53	2.03	1.95
14	A	602	HEA	CAA-C2A	2.50	1.57	1.51
14	A	602	HEA	CHC-C4B	2.50	1.43	1.38
23	G	104	CHD	C13-C17	-2.49	1.51	1.55
14	N	602	HEA	C3B-C2B	2.48	1.40	1.34
14	N	602	HEA	C3A-C4A	-2.47	1.41	1.46
19	A	607	TGL	OC1-CC1	-2.47	1.15	1.22
14	N	601	HEA	CHC-C4B	2.43	1.43	1.38
14	A	602	HEA	C4A-NA	-2.41	1.35	1.39
14	A	601	HEA	C4B-C3B	-2.41	1.40	1.44
14	A	601	HEA	C3B-C2B	2.38	1.40	1.34
14	N	602	HEA	C14-C15	2.37	1.38	1.33
14	N	602	HEA	O11-C11	2.34	1.48	1.42
20	C	302	PGV	O01-C1	2.30	1.40	1.34
14	A	601	HEA	C1D-C2D	-2.29	1.39	1.44
14	N	601	HEA	C3B-C2B	2.28	1.39	1.34
23	G	104	CHD	O7-C7	2.27	1.48	1.43
14	A	601	HEA	CHB-C1B	2.26	1.44	1.39
20	P	303	PGV	O03-C01	-2.24	1.40	1.45
14	N	601	HEA	C4A-NA	-2.23	1.35	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	A	601	HEA	C4D-ND	-2.22	1.34	1.38
14	N	601	HEA	O1D-CGD	2.22	1.29	1.22
14	N	601	HEA	C1D-C2D	-2.20	1.40	1.44
14	N	602	HEA	O1A-CGA	2.19	1.29	1.22
14	A	602	HEA	CHA-C4D	2.15	1.44	1.39
14	A	601	HEA	CHA-C4D	2.12	1.44	1.39
14	N	602	HEA	C4D-ND	-2.11	1.34	1.38
14	N	601	HEA	C1B-NB	-2.08	1.34	1.38
14	N	601	HEA	C4C-NC	-2.08	1.35	1.39
23	C	305	CHD	C4-C3	2.07	1.55	1.52
27	C	306	PEK	C01-C02	2.06	1.57	1.50
14	A	601	HEA	C1A-C2A	-2.04	1.41	1.45
14	N	601	HEA	C1A-C2A	-2.02	1.41	1.45
20	P	303	PGV	O01-C02	-2.02	1.41	1.46
14	A	602	HEA	C4B-C3B	-2.01	1.41	1.44

All (331) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	W	101	CHD	C14-C8-C9	-13.33	91.13	109.75
23	W	101	CHD	C18-C13-C12	-9.30	99.74	109.06
23	W	101	CHD	C14-C13-C12	9.29	115.91	107.42
19	A	607	TGL	OG2-CB1-CB2	8.20	129.22	111.48
23	W	101	CHD	C10-C9-C8	7.95	120.70	111.84
23	W	101	CHD	C17-C13-C14	-7.02	93.07	100.11
19	L	101	TGL	OG2-CB1-CB2	6.50	125.54	111.48
20	A	608	PGV	O03-C19-C20	6.17	130.64	111.83
23	W	101	CHD	C6-C7-C8	5.93	117.98	111.50
23	J	101	CHD	C13-C17-C20	5.89	126.62	119.48
14	A	602	HEA	C2A-C1A-NA	5.79	115.91	110.32
19	Y	101	TGL	OG2-CB1-CB2	5.67	123.75	111.48
26	P	304	CDL	OA6-CA5-C11	5.62	123.64	111.48
23	W	101	CHD	C13-C17-C20	5.58	126.24	119.48
19	A	607	TGL	CG3-CG2-CG1	-5.58	98.79	111.78
23	W	101	CHD	C14-C8-C7	5.43	119.06	111.85
20	A	608	PGV	O03-C19-O04	-5.43	110.05	123.63
23	W	101	CHD	C17-C13-C12	5.36	122.49	117.67
23	J	101	CHD	C17-C13-C14	-5.26	94.84	100.11
19	O	302	TGL	OG2-CB1-CB2	5.24	122.83	111.48
14	N	601	HEA	C3C-C4C-NC	5.24	114.21	109.80
14	A	602	HEA	CMD-C2D-C1D	5.23	133.20	125.03
23	W	101	CHD	C9-C11-C12	-5.20	107.48	114.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	A	602	HEA	C27-C19-C20	5.19	124.24	115.23
23	W	101	CHD	C15-C14-C8	5.11	125.37	118.36
20	P	301	PGV	O01-C1-C2	5.09	122.49	111.48
14	A	602	HEA	C3A-C2A-C1A	-5.08	102.24	107.05
23	P	306	CHD	C15-C14-C13	5.01	108.40	103.54
26	G	102	CDL	OB6-CB5-C51	5.00	122.29	111.48
20	N	607	PGV	O01-C1-C2	4.97	122.23	111.48
27	G	101	PEK	O01-C1-C2	4.92	122.13	111.48
26	C	303	CDL	OA6-CA5-C11	4.84	121.94	111.48
23	W	101	CHD	C9-C8-C7	4.77	117.86	111.86
14	N	602	HEA	CMB-C2B-C1B	4.72	132.41	125.03
23	W	101	CHD	C11-C9-C10	4.67	118.44	113.70
26	T	104	CDL	OB6-CB5-C51	4.52	121.27	111.48
27	C	306	PEK	O01-C1-C2	4.49	121.18	111.48
23	W	101	CHD	C11-C12-C13	4.46	115.80	111.26
26	P	304	CDL	OB6-CB5-C51	4.44	121.08	111.48
19	L	101	TGL	OG3-CC1-CC2	4.42	125.30	111.83
24	B	303	PSC	O01-C1-C2	4.33	120.86	111.48
19	A	607	TGL	OG3-CC1-OC1	-4.30	112.87	123.63
27	G	103	PEK	O01-C1-C2	4.28	120.74	111.48
23	C	304	CHD	C15-C14-C13	4.27	107.68	103.54
14	N	602	HEA	CAD-CBD-CGD	-4.22	102.47	113.67
23	P	305	CHD	C14-C13-C12	4.20	111.25	107.42
27	T	103	PEK	O01-C1-C2	4.19	120.56	111.48
23	C	305	CHD	C21-C20-C22	-4.18	103.88	110.34
23	C	305	CHD	C23-C22-C20	-4.16	106.69	114.46
14	A	601	HEA	C13-C12-C11	-4.16	107.75	114.39
14	N	602	HEA	CHB-C4A-NA	4.13	128.95	124.45
23	G	104	CHD	C4-C5-C10	-4.08	108.32	112.66
27	T	102	PEK	O01-C1-C2	4.02	120.18	111.48
20	A	609	PGV	O01-C1-C2	4.00	120.13	111.48
26	C	303	CDL	OB6-CB5-C51	3.99	120.11	111.48
23	P	305	CHD	C9-C10-C5	3.99	114.05	108.51
27	G	101	PEK	C01-O03-C21	3.99	131.70	117.12
19	O	302	TGL	OG3-CC1-CC2	3.97	123.94	111.83
27	T	101	PEK	O03-C01-C02	-3.93	97.07	108.40
26	G	102	CDL	OA6-CA5-C11	3.93	119.97	111.48
19	A	607	TGL	OG3-CC1-CC2	3.92	123.79	111.83
24	O	303	PSC	O03-C19-C20	3.91	123.74	111.83
14	N	602	HEA	CAD-C3D-C2D	3.90	135.17	127.87
14	N	602	HEA	O11-C11-C3B	-3.88	104.14	111.26
14	N	602	HEA	CHC-C4B-NB	3.87	129.16	124.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
19	D	201	TGL	OG3-CC1-CC2	3.85	123.57	111.83
14	A	602	HEA	CHD-C4C-NC	3.83	130.81	123.86
23	W	101	CHD	C16-C17-C13	3.83	107.26	103.54
26	T	104	CDL	OA6-CA5-C11	3.82	119.75	111.48
23	J	101	CHD	C17-C13-C12	3.82	121.10	117.67
20	A	609	PGV	O03-C19-C20	3.81	123.46	111.83
19	Y	101	TGL	OG3-CC1-CC2	3.81	123.45	111.83
26	C	303	CDL	OB8-CB7-C71	3.78	123.35	111.83
14	A	601	HEA	C26-C15-C16	3.77	121.78	115.23
14	A	601	HEA	CAA-C2A-C1A	3.77	132.52	124.85
23	B	302	CHD	C15-C14-C13	3.76	107.18	103.54
14	N	602	HEA	C25-C23-C24	3.74	123.20	114.59
19	Q	201	TGL	OG2-CB1-CB2	3.74	119.57	111.48
14	A	602	HEA	CAD-CBD-CGD	-3.72	103.81	113.67
14	A	601	HEA	C21-C22-C23	3.70	139.98	127.64
26	P	304	CDL	OB8-CB7-C71	3.68	123.04	111.83
14	A	602	HEA	CHA-C1A-C2A	-3.64	119.11	124.86
27	G	101	PEK	O01-C1-O02	-3.64	115.20	123.70
14	A	602	HEA	C26-C15-C16	3.64	121.54	115.23
23	P	306	CHD	C17-C13-C14	-3.63	96.48	100.11
14	A	601	HEA	C21-C20-C19	-3.61	101.23	113.19
23	W	101	CHD	C22-C23-C24	-3.60	102.90	112.49
14	N	602	HEA	C1C-CHC-C4B	-3.57	117.86	126.25
26	T	104	CDL	OB8-CB7-C71	3.55	122.66	111.83
27	C	306	PEK	O03-C01-C02	3.47	118.40	108.40
23	C	304	CHD	C13-C14-C8	-3.45	110.34	114.72
20	N	608	PGV	O03-C19-C20	3.44	122.33	111.83
20	P	301	PGV	O03-C19-C20	3.44	122.31	111.83
23	C	304	CHD	C6-C5-C4	-3.43	107.31	111.23
23	C	304	CHD	C1-C10-C5	3.41	112.64	107.75
23	P	305	CHD	C1-C10-C9	-3.39	106.08	111.34
26	G	102	CDL	OB8-CB7-C71	3.38	122.13	111.83
24	O	303	PSC	O01-C1-C2	3.37	118.78	111.48
19	O	302	TGL	OG1-CA1-CA2	3.37	122.10	111.83
23	C	304	CHD	C10-C9-C8	3.36	115.58	111.84
20	C	307	PGV	O01-C1-C2	3.35	118.74	111.48
14	A	602	HEA	C4A-NA-C1A	-3.35	100.35	105.82
23	P	305	CHD	C15-C14-C13	3.33	106.77	103.54
26	G	102	CDL	OA8-CA7-C31	3.32	121.95	111.83
24	B	303	PSC	O03-C19-C20	3.31	121.92	111.83
14	A	602	HEA	C1D-C2D-C3D	-3.31	103.50	106.98
19	A	607	TGL	OG1-CA1-CA2	3.28	121.85	111.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
26	P	304	CDL	OA8-CA7-C31	3.27	121.81	111.83
20	N	608	PGV	O01-C1-C2	3.27	118.56	111.48
14	A	602	HEA	C4C-CHD-C1D	-3.26	118.58	126.25
19	L	101	TGL	CG2-OG2-CB1	3.26	125.60	117.80
23	C	304	CHD	C11-C9-C10	-3.26	110.40	113.70
23	W	101	CHD	C1-C10-C9	-3.25	106.29	111.34
26	P	304	CDL	OB8-CB7-OB9	-3.23	115.55	123.63
14	N	602	HEA	CMB-C2B-C3B	-3.21	124.06	130.28
14	A	602	HEA	C3C-C4C-NC	3.21	112.50	109.80
26	C	303	CDL	OA8-CA7-C31	3.19	121.55	111.83
14	A	602	HEA	C13-C12-C11	-3.16	109.35	114.39
20	P	303	PGV	C03-C02-C01	-3.14	104.45	111.78
20	N	608	PGV	O03-C19-O04	-3.14	115.78	123.63
23	C	304	CHD	C6-C5-C10	3.13	115.99	112.66
23	B	302	CHD	C6-C5-C4	-3.09	107.69	111.23
23	P	305	CHD	C6-C5-C10	3.08	115.94	112.66
19	D	201	TGL	CG1-OG1-CA1	3.07	128.35	117.12
19	Q	201	TGL	OG3-CC1-CC2	3.06	121.17	111.83
19	Y	101	TGL	OG3-CC1-OC1	-3.06	115.98	123.63
14	A	602	HEA	CMB-C2B-C1B	3.04	129.79	125.03
27	C	306	PEK	O03-C21-C22	3.02	121.06	111.83
23	P	306	CHD	C5-C6-C7	-3.02	110.81	114.40
14	A	601	HEA	C26-C15-C14	-3.01	115.89	123.63
20	P	303	PGV	O01-C1-C2	3.01	118.00	111.48
23	W	101	CHD	C16-C15-C14	-3.00	99.27	105.14
19	Q	201	TGL	OG1-CA1-CA2	2.99	120.97	111.83
19	A	607	TGL	OB1-CB1-CB2	-2.98	112.14	123.78
23	P	306	CHD	C1-C10-C5	2.97	112.01	107.75
23	C	304	CHD	C9-C10-C5	2.97	112.64	108.51
14	N	602	HEA	O11-C11-C12	2.96	116.98	109.14
23	J	101	CHD	C22-C20-C17	2.95	116.45	110.33
14	N	601	HEA	C1D-C2D-C3D	-2.94	103.89	106.98
23	C	304	CHD	C4-C5-C10	2.94	115.79	112.66
14	A	601	HEA	C3C-C4C-NC	2.93	112.27	109.80
14	A	601	HEA	CHA-C4D-C3D	-2.93	120.50	124.77
23	G	104	CHD	C13-C17-C20	-2.93	115.94	119.48
20	C	307	PGV	O03-C19-C20	2.92	120.75	111.83
27	T	102	PEK	O03-C21-C22	2.92	120.75	111.83
20	A	609	PGV	O03-C19-O04	-2.92	116.33	123.63
23	C	304	CHD	C2-C1-C10	2.90	117.64	112.74
29	M	101	DMU	O1-C9-C11	2.90	113.63	106.44
19	Q	201	TGL	CG2-OG2-CB1	-2.90	110.87	117.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	C	305	CHD	C6-C5-C10	2.89	115.73	112.66
14	A	602	HEA	C1C-CHC-C4B	-2.88	119.46	126.25
14	A	602	HEA	C27-C19-C18	-2.88	116.23	123.63
23	C	304	CHD	C15-C14-C8	2.88	122.31	118.36
14	N	602	HEA	CHA-C1A-C2A	-2.87	120.33	124.86
23	C	305	CHD	O25-C24-C23	-2.87	113.99	123.09
20	N	608	PGV	O01-C1-O02	-2.86	117.01	123.70
20	P	303	PGV	O03-C19-O04	-2.85	116.49	123.63
14	N	601	HEA	C17-C18-C19	-2.85	121.10	127.62
14	A	602	HEA	CMC-C2C-C1C	2.85	129.75	125.42
14	N	602	HEA	C20-C19-C18	-2.85	114.78	121.17
14	N	602	HEA	C1B-C2B-C3B	-2.83	103.52	106.80
14	A	602	HEA	CHC-C4B-NB	2.82	127.85	124.37
27	T	103	PEK	O03-C21-C22	2.81	120.42	111.83
23	P	305	CHD	C10-C9-C8	2.81	114.97	111.84
23	P	306	CHD	C18-C13-C14	2.81	115.61	111.20
23	G	104	CHD	O12-C12-C13	-2.81	106.28	111.02
23	P	305	CHD	C11-C9-C10	-2.81	110.85	113.70
23	C	304	CHD	C16-C17-C20	2.81	116.42	112.18
27	G	103	PEK	O03-C21-C22	2.80	120.36	111.83
23	C	305	CHD	O26-C24-C23	2.80	122.83	114.00
19	D	201	TGL	OG2-CB1-CB2	2.79	117.52	111.48
23	W	101	CHD	C23-C22-C20	-2.79	109.25	114.46
23	C	305	CHD	C19-C10-C1	-2.79	103.87	108.31
19	O	302	TGL	CG3-CG2-CG1	-2.79	105.29	111.78
23	G	104	CHD	O26-C24-C23	2.77	122.77	114.00
19	A	607	TGL	OG2-CG2-CG3	2.77	118.29	108.34
14	N	601	HEA	C3A-C2A-C1A	-2.77	104.42	107.05
14	N	602	HEA	C1A-CHA-C4D	-2.77	120.13	126.02
27	G	103	PEK	O03-C01-C02	2.77	116.38	108.40
23	W	101	CHD	C5-C6-C7	2.76	117.68	114.40
20	P	303	PGV	C04-C05-C06	-2.75	102.51	111.77
26	T	104	CDL	OA8-CA7-C31	2.75	120.22	111.83
14	N	601	HEA	CMC-C2C-C1C	2.75	129.60	125.42
26	G	102	CDL	CB4-OB6-CB5	2.73	124.34	117.80
19	Q	201	TGL	OG3-CC1-OC1	-2.73	116.81	123.63
20	C	302	PGV	O03-C19-O04	-2.71	116.84	123.63
14	A	602	HEA	CBD-CAD-C3D	2.70	120.00	112.53
14	N	602	HEA	CHC-C1C-C2C	-2.70	119.58	127.43
14	N	602	HEA	CHB-C4A-C3A	-2.69	120.68	125.21
19	D	201	TGL	CG3-OG3-CC1	2.67	126.89	117.12
14	A	602	HEA	CHD-C4C-C3C	-2.67	115.80	127.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
23	W	101	CHD	C21-C20-C17	2.65	116.87	112.88
14	N	602	HEA	C25-C23-C22	-2.65	114.69	122.66
23	J	101	CHD	C6-C7-C8	2.64	114.39	111.50
20	P	303	PGV	O03-C19-C20	2.64	119.87	111.83
23	B	302	CHD	O26-C24-C23	2.63	122.30	114.00
27	G	103	PEK	O01-C1-O02	-2.62	117.57	123.70
19	Y	101	TGL	OG1-CA1-CA2	2.62	119.83	111.83
23	W	101	CHD	C4-C5-C10	2.62	115.44	112.66
19	Y	101	TGL	OG2-CB1-OB1	-2.62	117.59	123.70
24	B	303	PSC	O01-C1-O02	-2.59	117.65	123.70
23	J	101	CHD	C1-C10-C5	2.59	111.46	107.75
23	B	302	CHD	C19-C10-C5	-2.58	106.13	110.44
20	N	607	PGV	O03-C01-C02	2.56	115.78	108.40
26	P	304	CDL	OB6-CB5-OB7	-2.56	117.72	123.70
19	L	101	TGL	OG2-CB1-OB1	-2.53	117.78	123.70
14	N	602	HEA	CAD-C3D-C4D	-2.53	120.29	124.70
19	D	201	TGL	OG3-CC1-OC1	-2.52	117.32	123.63
14	N	601	HEA	C2D-C1D-ND	2.52	112.73	109.84
26	C	303	CDL	OB8-CB7-OB9	-2.51	117.35	123.63
23	P	305	CHD	C1-C10-C5	2.50	111.34	107.75
23	B	302	CHD	C6-C7-C8	2.49	114.22	111.50
23	G	104	CHD	O3-C3-C4	-2.48	104.90	109.84
20	P	303	PGV	O14-P-O13	2.47	123.94	112.44
23	C	304	CHD	C16-C17-C13	2.47	105.93	103.54
23	C	305	CHD	C15-C14-C13	2.45	105.92	103.54
23	P	306	CHD	C19-C10-C5	-2.45	106.35	110.44
20	A	608	PGV	O03-C01-C02	2.43	115.41	108.40
27	T	101	PEK	O01-C1-C2	2.43	116.73	111.48
27	T	101	PEK	C01-O03-C21	2.43	125.99	117.12
23	G	104	CHD	C13-C14-C8	-2.42	111.65	114.72
20	A	609	PGV	C02-O01-C1	2.42	123.59	117.80
19	D	201	TGL	OG1-CA1-CA2	2.42	119.21	111.83
23	J	101	CHD	C6-C5-C10	2.42	115.23	112.66
19	L	101	TGL	CA4-CA3-CA2	-2.41	104.29	113.13
14	A	602	HEA	C25-C23-C24	2.40	120.12	114.59
14	N	601	HEA	CHC-C4B-NB	2.40	127.33	124.37
19	O	302	TGL	OG2-CB1-OB1	-2.40	118.10	123.70
26	C	303	CDL	OA6-CA5-OA7	-2.39	118.12	123.70
20	P	301	PGV	C01-O03-C19	2.39	125.85	117.12
26	G	102	CDL	CB6-OB8-CB7	2.39	125.85	117.12
23	B	302	CHD	C11-C9-C10	2.39	116.13	113.70
19	O	302	TGL	OG3-CG3-CG2	2.38	115.26	108.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
24	O	303	PSC	O03-C01-C02	2.38	115.26	108.40
23	B	302	CHD	C13-C14-C8	-2.37	111.71	114.72
14	N	602	HEA	C2B-C1B-NB	2.37	112.65	109.90
23	P	306	CHD	C19-C10-C1	-2.37	104.54	108.31
20	C	302	PGV	O03-C19-C20	2.37	119.06	111.83
14	A	602	HEA	CAD-C3D-C4D	2.37	128.83	124.70
23	P	305	CHD	C14-C8-C9	-2.37	106.44	109.75
14	N	602	HEA	C4A-CHB-C1B	-2.37	120.99	126.02
14	N	601	HEA	C4C-C3C-C2C	-2.36	104.37	107.30
23	G	104	CHD	C2-C1-C10	2.36	116.72	112.74
23	P	306	CHD	C11-C9-C8	2.36	114.38	110.89
20	P	301	PGV	O01-C1-O02	-2.35	118.20	123.70
19	L	101	TGL	OG1-CA1-CA2	2.35	119.01	111.83
23	G	104	CHD	O7-C7-C6	2.35	115.71	109.86
14	N	601	HEA	CAA-C2A-C1A	2.35	129.63	124.85
14	A	602	HEA	CMB-C2B-C3B	-2.35	125.73	130.28
14	A	601	HEA	CHA-C4D-ND	2.35	126.95	124.42
23	J	101	CHD	C14-C13-C12	2.34	109.55	107.42
23	P	305	CHD	C4-C3-C2	2.33	113.46	110.62
14	A	601	HEA	CHB-C4A-NA	2.33	126.99	124.45
23	P	305	CHD	C18-C13-C12	-2.33	106.72	109.06
14	N	601	HEA	C4B-NB-C1B	2.33	107.96	105.21
26	G	102	CDL	OB8-CB6-CB4	2.32	115.09	108.40
14	N	602	HEA	C13-C12-C11	-2.31	110.70	114.39
19	Y	101	TGL	OG3-CG3-CG2	2.31	115.05	108.40
14	N	602	HEA	CAA-C2A-C1A	2.30	129.54	124.85
14	A	602	HEA	C2D-C1D-ND	2.30	112.49	109.84
23	B	302	CHD	C13-C17-C20	-2.30	116.69	119.48
19	L	101	TGL	OG3-CG3-CG2	2.30	115.01	108.40
26	G	102	CDL	OA8-CA7-OA9	-2.29	117.90	123.63
20	N	607	PGV	O03-C19-C20	2.29	118.82	111.83
14	N	602	HEA	CHC-C4B-C3B	-2.29	120.02	125.80
23	P	306	CHD	C5-C4-C3	-2.29	109.26	112.71
20	P	303	PGV	C22-C21-C20	-2.28	104.76	113.13
20	A	609	PGV	C6-C5-C4	-2.27	102.89	114.37
14	A	601	HEA	C12-C11-C3B	2.27	115.67	112.12
14	N	601	HEA	CHB-C4A-NA	2.26	126.92	124.45
20	C	307	PGV	C01-O03-C19	2.26	125.38	117.12
27	T	102	PEK	O01-C1-O02	-2.26	118.42	123.70
23	G	104	CHD	C15-C14-C8	-2.26	115.26	118.36
23	C	304	CHD	C9-C8-C7	2.25	114.69	111.86
23	P	306	CHD	C14-C8-C7	2.25	114.83	111.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	A	601	HEA	CHD-C4C-NC	2.24	127.93	123.86
26	T	104	CDL	CA6-OA8-CA7	2.24	125.30	117.12
20	N	607	PGV	C03-C02-C01	-2.24	106.57	111.78
26	C	303	CDL	CB6-OB8-CB7	2.23	125.27	117.12
26	G	102	CDL	OB6-CB5-OB7	-2.22	118.51	123.70
23	B	302	CHD	C11-C9-C8	2.22	114.18	110.89
23	B	302	CHD	C4-C3-C2	-2.21	107.92	110.62
27	T	102	PEK	O03-C21-O04	-2.21	118.10	123.63
23	P	305	CHD	C6-C5-C4	-2.21	108.70	111.23
26	P	304	CDL	OA8-CA7-OA9	-2.21	118.11	123.63
14	A	602	HEA	O2A-CGA-CBA	2.20	120.96	114.00
23	P	306	CHD	C6-C5-C4	2.20	113.74	111.23
23	J	101	CHD	C13-C14-C8	2.20	117.50	114.72
27	G	101	PEK	C3-C2-C1	-2.19	105.66	113.69
27	C	306	PEK	O01-C02-C01	2.19	116.21	108.34
26	C	303	CDL	OB8-CB6-CB4	2.19	114.70	108.40
14	A	601	HEA	CAA-CBA-CGA	-2.19	107.87	113.67
19	L	101	TGL	OG3-CC1-OC1	-2.18	118.17	123.63
19	A	607	TGL	OG2-CB1-OB1	-2.17	118.62	123.70
24	O	303	PSC	O03-C19-O04	-2.17	118.19	123.63
23	W	101	CHD	C22-C20-C17	2.17	114.83	110.33
27	C	306	PEK	O03-C21-O04	-2.16	118.23	123.63
23	W	101	CHD	O3-C3-C4	2.16	114.14	109.84
27	G	103	PEK	O01-C02-C01	2.15	116.07	108.34
14	N	602	HEA	CHA-C1A-NA	2.15	126.79	124.45
14	A	601	HEA	O2D-CGD-O1D	-2.14	117.83	123.33
14	N	602	HEA	CHA-C4D-C3D	-2.14	121.66	124.77
19	O	302	TGL	OG3-CC1-OC1	-2.13	118.29	123.63
23	C	305	CHD	C17-C13-C12	2.13	119.58	117.67
29	Z	101	DMU	O49-C1-C6	-2.12	105.02	110.08
23	C	304	CHD	C19-C10-C9	-2.12	108.33	111.18
20	C	302	PGV	O01-C1-O02	-2.11	118.76	123.70
20	N	607	PGV	O01-C1-O02	-2.11	118.78	123.70
23	P	305	CHD	C15-C14-C8	2.10	121.24	118.36
14	N	601	HEA	CAA-CBA-CGA	-2.10	108.10	113.67
14	A	602	HEA	C1B-C2B-C3B	-2.10	104.36	106.80
23	W	101	CHD	C18-C13-C17	2.10	114.49	111.20
26	C	303	CDL	OA8-CA7-OA9	-2.10	118.38	123.63
23	C	305	CHD	C14-C13-C12	-2.09	105.50	107.42
20	A	609	PGV	O01-C02-C03	2.09	115.83	108.34
23	C	304	CHD	C19-C10-C1	-2.09	104.99	108.31
24	B	303	PSC	O01-C02-C01	2.08	115.82	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	A	602	HEA	C21-C20-C19	2.08	120.09	113.19
14	A	601	HEA	C3D-C4D-ND	2.08	112.36	110.35
23	C	305	CHD	C1-C10-C5	2.07	110.72	107.75
29	Z	101	DMU	O1-C9-C11	2.06	111.56	106.44
23	B	302	CHD	C16-C17-C13	2.06	105.54	103.54
23	C	305	CHD	C18-C13-C14	2.06	114.43	111.20
27	T	102	PEK	C01-O03-C21	2.04	124.59	117.12
20	C	302	PGV	O01-C1-C2	2.04	115.89	111.48
14	A	601	HEA	C3A-C2A-C1A	-2.03	105.13	107.05
23	C	304	CHD	O25-C24-C23	-2.02	116.67	123.09
27	G	101	PEK	C02-O01-C1	-2.02	112.96	117.80
23	G	104	CHD	C11-C9-C8	2.02	113.88	110.89
20	A	608	PGV	C5-C4-C3	-2.01	104.22	114.37
23	G	104	CHD	C18-C13-C12	-2.00	107.05	109.06

There are no chirality outliers.

All (941) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	N	601	HEA	C2A-C3A-CMA-OMA
19	D	201	TGL	CB2-CB1-OG2-CG2
19	D	201	TGL	OB1-CB1-OG2-CG2
19	L	101	TGL	CB2-CB1-OG2-CG2
19	L	101	TGL	OB1-CB1-OG2-CG2
19	Q	201	TGL	CC2-CC1-OG3-CG3
19	Q	201	TGL	OC1-CC1-OG3-CG3
19	Y	101	TGL	CB2-CB1-OG2-CG2
19	Y	101	TGL	OB1-CB1-OG2-CG2
20	A	609	PGV	C04-O12-P-O11
20	A	609	PGV	C02-C03-O11-P
20	A	609	PGV	O02-C1-O01-C02
20	A	609	PGV	C2-C1-O01-C02
20	C	307	PGV	C03-O11-P-O12
20	C	307	PGV	C03-O11-P-O13
20	C	307	PGV	C03-O11-P-O14
20	C	307	PGV	O12-C04-C05-C06
20	C	307	PGV	C04-C05-C06-O06
20	N	607	PGV	C04-O12-P-O13
20	N	607	PGV	C02-C03-O11-P
20	N	607	PGV	O12-C04-C05-C06
20	N	607	PGV	C2-C1-O01-C02
20	P	301	PGV	C03-O11-P-O12

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Mol	Chain	Res	Type	Atoms
20	P	301	PGV	C2-C1-O01-C02
23	W	101	CHD	C13-C17-C20-C22
24	B	303	PSC	C03-O11-P-O12
24	B	303	PSC	C04-O12-P-O14
24	B	303	PSC	C2-C1-O01-C02
24	O	303	PSC	C03-O11-P-O12
24	O	303	PSC	C03-O11-P-O13
24	O	303	PSC	C04-O12-P-O11
24	O	303	PSC	C04-O12-P-O13
24	O	303	PSC	C04-O12-P-O14
24	O	303	PSC	O12-C04-C05-N
24	O	303	PSC	C11-C12-C13-C14
26	C	303	CDL	CA2-C1-CB2-OB2
26	C	303	CDL	CA2-OA2-PA1-OA3
26	C	303	CDL	CA2-OA2-PA1-OA4
26	C	303	CDL	CA2-OA2-PA1-OA5
26	C	303	CDL	CA3-OA5-PA1-OA2
26	C	303	CDL	C11-CA5-OA6-CA4
26	C	303	CDL	C51-CB5-OB6-CB4
26	G	102	CDL	O1-C1-CA2-OA2
26	G	102	CDL	CA2-OA2-PA1-OA3
26	G	102	CDL	CA2-OA2-PA1-OA4
26	G	102	CDL	CA2-OA2-PA1-OA5
26	G	102	CDL	C11-CA5-OA6-CA4
26	G	102	CDL	CB3-OB5-PB2-OB2
26	G	102	CDL	OB7-CB5-OB6-CB4
26	G	102	CDL	C51-CB5-OB6-CB4
26	P	304	CDL	CA2-OA2-PA1-OA3
26	P	304	CDL	CA2-OA2-PA1-OA4
26	P	304	CDL	CA2-OA2-PA1-OA5
26	P	304	CDL	CA3-OA5-PA1-OA2
26	P	304	CDL	CA3-OA5-PA1-OA3
26	P	304	CDL	CA3-OA5-PA1-OA4
26	P	304	CDL	OA7-CA5-OA6-CA4
26	P	304	CDL	C11-CA5-OA6-CA4
26	P	304	CDL	CB3-OB5-PB2-OB2
26	P	304	CDL	CB3-OB5-PB2-OB4
26	T	104	CDL	CA2-C1-CB2-OB2
26	T	104	CDL	CA2-OA2-PA1-OA3
26	T	104	CDL	CA2-OA2-PA1-OA4
26	T	104	CDL	C11-CA5-OA6-CA4
26	T	104	CDL	CB3-OB5-PB2-OB2

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Mol	Chain	Res	Type	Atoms
26	T	104	CDL	CB3-OB5-PB2-OB3
27	C	306	PEK	C03-O11-P-O14
27	C	306	PEK	C04-O12-P-O14
27	C	306	PEK	O12-C04-C05-N
27	G	103	PEK	C04-O12-P-O11
27	G	103	PEK	C04-O12-P-O13
27	G	103	PEK	C04-O12-P-O14
27	G	103	PEK	O12-C04-C05-N
27	G	103	PEK	C2-C1-O01-C02
27	T	102	PEK	C03-O11-P-O12
27	T	102	PEK	C03-O11-P-O13
27	T	102	PEK	C12-C13-C14-C15
27	T	103	PEK	C04-O12-P-O11
27	T	103	PEK	C04-O12-P-O13
27	T	103	PEK	O03-C01-C02-O01
27	T	103	PEK	C5-C6-C7-C8
27	T	103	PEK	C9-C10-C11-C12
19	D	201	TGL	OC1-CC1-OG3-CG3
24	O	303	PSC	O04-C19-O03-C01
24	O	303	PSC	C20-C19-O03-C01
19	O	302	TGL	OC1-CC1-OG3-CG3
20	A	609	PGV	O04-C19-O03-C01
20	N	607	PGV	O04-C19-O03-C01
24	B	303	PSC	O04-C19-O03-C01
26	T	104	CDL	OA9-CA7-OA8-CA6
27	T	102	PEK	O04-C21-O03-C01
20	N	607	PGV	O02-C1-O01-C02
20	P	301	PGV	O02-C1-O01-C02
26	C	303	CDL	OA7-CA5-OA6-CA4
26	C	303	CDL	OB7-CB5-OB6-CB4
27	G	103	PEK	O02-C1-O01-C02
19	A	607	TGL	CC2-CC1-OG3-CG3
19	D	201	TGL	CC2-CC1-OG3-CG3
19	O	302	TGL	CC2-CC1-OG3-CG3
20	A	609	PGV	C20-C19-O03-C01
26	T	104	CDL	C31-CA7-OA8-CA6
19	A	607	TGL	CA2-CA1-OG1-CG1
20	N	607	PGV	C20-C19-O03-C01
24	B	303	PSC	C20-C19-O03-C01
26	C	303	CDL	C71-CB7-OB8-CB6
27	T	102	PEK	C22-C21-O03-C01
27	G	101	PEK	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
27	G	101	PEK	C7-C8-C9-C10
26	C	303	CDL	OB9-CB7-OB8-CB6
24	B	303	PSC	O02-C1-O01-C02
26	G	102	CDL	OA7-CA5-OA6-CA4
26	T	104	CDL	OA7-CA5-OA6-CA4
24	B	303	PSC	C04-C05-N-C08
20	N	607	PGV	O12-C04-C05-O05
26	G	102	CDL	O1-C1-CB2-OB2
26	P	304	CDL	O1-C1-CA2-OA2
19	A	607	TGL	OC1-CC1-OG3-CG3
19	Q	201	TGL	CB2-CB1-OG2-CG2
19	A	607	TGL	OA1-CA1-OG1-CG1
23	J	101	CHD	C17-C20-C22-C23
23	J	101	CHD	C21-C20-C22-C23
23	P	305	CHD	C21-C20-C22-C23
24	O	303	PSC	C24-C25-C26-C27
19	L	101	TGL	CA2-CA1-OG1-CG1
23	P	305	CHD	C17-C20-C22-C23
20	A	609	PGV	C10-C11-C12-C13
20	C	307	PGV	C10-C11-C12-C13
27	G	101	PEK	C13-C14-C15-C16
27	T	101	PEK	C13-C14-C15-C16
19	A	607	TGL	CB2-CB1-OG2-CG2
20	A	609	PGV	O12-C04-C05-C06
26	C	303	CDL	CB2-C1-CA2-OA2
26	G	102	CDL	CB2-C1-CA2-OA2
26	P	304	CDL	CB2-C1-CA2-OA2
26	G	102	CDL	C31-CA7-OA8-CA6
19	O	302	TGL	CA7-CA8-CA9-C20
29	M	101	DMU	O6-C11-C9-O1
20	A	608	PGV	C26-C27-C28-C29
26	G	102	CDL	C78-C79-C80-C81
19	Y	101	TGL	CA9-C20-C21-C22
29	M	101	DMU	O6-C11-C9-C8
20	C	307	PGV	O12-C04-C05-O05
26	C	303	CDL	O1-C1-CB2-OB2
26	T	104	CDL	O1-C1-CB2-OB2
20	A	609	PGV	C20-C21-C22-C23
26	G	102	CDL	OA9-CA7-OA8-CA6
19	Q	201	TGL	OB1-CB1-OG2-CG2
26	P	304	CDL	C51-CB5-OB6-CB4
26	C	303	CDL	CA5-C11-C12-C13

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Mol	Chain	Res	Type	Atoms
27	C	306	PEK	C4-C5-C6-C7
27	C	306	PEK	C10-C11-C12-C13
20	A	609	PGV	O05-C05-C06-O06
19	O	302	TGL	CA2-CA1-OG1-CG1
27	G	101	PEK	C22-C21-O03-C01
19	L	101	TGL	CB1-CB2-CB3-CB4
19	D	201	TGL	CB1-CB2-CB3-CB4
19	Q	201	TGL	CC1-CC2-CC3-CC4
20	P	303	PGV	C1-C2-C3-C4
24	B	303	PSC	C1-C2-C3-C4
26	P	304	CDL	CB7-C71-C72-C73
26	P	304	CDL	C1-CA2-OA2-PA1
19	Y	101	TGL	CA1-CA2-CA3-CA4
20	N	607	PGV	C19-C20-C21-C22
27	G	103	PEK	C21-C22-C23-C24
19	L	101	TGL	OA1-CA1-OG1-CG1
27	G	101	PEK	O04-C21-O03-C01
26	C	303	CDL	O1-C1-CA2-OA2
26	P	304	CDL	OB7-CB5-OB6-CB4
20	N	608	PGV	C10-C11-C12-C13
20	P	303	PGV	C10-C11-C12-C13
24	B	303	PSC	C11-C10-C9-C8
27	T	103	PEK	C13-C14-C15-C16
19	D	201	TGL	CA1-CA2-CA3-CA4
26	T	104	CDL	CA7-C31-C32-C33
26	T	104	CDL	C71-CB7-OB8-CB6
26	C	303	CDL	C59-C60-C61-C62
19	Y	101	TGL	CB1-CB2-CB3-CB4
19	A	607	TGL	OB1-CB1-OG2-CG2
26	G	102	CDL	CA2-C1-CB2-OB2
19	Y	101	TGL	CA2-CA1-OG1-CG1
23	W	101	CHD	C16-C17-C20-C21
24	B	303	PSC	C04-C05-N-C06
19	L	101	TGL	CC1-CC2-CC3-CC4
20	A	608	PGV	C10-C11-C12-C13
27	G	103	PEK	C4-C5-C6-C7
27	G	103	PEK	C10-C11-C12-C13
20	A	609	PGV	O12-C04-C05-O05
19	O	302	TGL	OA1-CA1-OG1-CG1
20	A	609	PGV	C04-C05-C06-O06
20	N	607	PGV	C03-C02-O01-C1
27	G	103	PEK	C01-C02-O01-C1

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Mol	Chain	Res	Type	Atoms
26	T	104	CDL	OB9-CB7-OB8-CB6
26	T	104	CDL	OB7-CB5-OB6-CB4
27	C	306	PEK	O03-C01-C02-O01
27	C	306	PEK	C22-C21-O03-C01
24	B	303	PSC	C04-C05-N-C07
29	M	101	DMU	O16-C18-C19-C22
23	J	101	CHD	C13-C17-C20-C22
19	Q	201	TGL	C11-C10-CB9-CB8
19	Y	101	TGL	C23-C24-C25-C26
19	Y	101	TGL	CB9-C10-C11-C12
20	A	609	PGV	C14-C15-C16-C17
20	C	307	PGV	C24-C25-C26-C27
24	B	303	PSC	C4-C5-C6-C7
24	B	303	PSC	C26-C27-C28-C29
26	P	304	CDL	C60-C61-C62-C63
29	Z	101	DMU	C19-C22-C25-C28
27	G	103	PEK	C7-C8-C9-C10
27	T	101	PEK	C10-C11-C12-C13
19	L	101	TGL	CB4-CB5-CB6-CB7
19	Y	101	TGL	C11-C12-C13-C14
27	T	103	PEK	C34-C35-C36-C37
29	Z	101	DMU	C31-C34-C37-C40
19	A	607	TGL	C20-C21-C22-C23
19	Q	201	TGL	CA6-CA7-CA8-CA9
19	Y	101	TGL	CB6-CB7-CB8-CB9
26	G	102	CDL	C55-C56-C57-C58
26	G	102	CDL	C77-C78-C79-C80
26	T	104	CDL	C32-C33-C34-C35
26	T	104	CDL	C79-C80-C81-C82
19	Y	101	TGL	CB3-CB4-CB5-CB6
26	C	303	CDL	C80-C81-C82-C83
20	C	302	PGV	C1-C2-C3-C4
26	G	102	CDL	CA5-C11-C12-C13
19	D	201	TGL	CB3-CB4-CB5-CB6
19	Y	101	TGL	C21-C22-C23-C24
20	A	608	PGV	C4-C5-C6-C7
23	W	101	CHD	C16-C17-C20-C22
24	O	303	PSC	C20-C21-C22-C23
26	T	104	CDL	C19-C20-C21-C22
19	O	302	TGL	CB2-CB1-OG2-CG2
26	T	104	CDL	C51-CB5-OB6-CB4
26	T	104	CDL	C53-C54-C55-C56

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Mol	Chain	Res	Type	Atoms
23	W	101	CHD	C17-C20-C22-C23
19	D	201	TGL	C18-C19-C33-C34
19	L	101	TGL	CB5-CB6-CB7-CB8
20	P	301	PGV	C26-C27-C28-C29
27	G	103	PEK	C30-C31-C32-C33
19	L	101	TGL	CB7-CB8-CB9-C10
20	P	303	PGV	C7-C8-C9-C10
26	G	102	CDL	C33-C34-C35-C36
26	G	102	CDL	C59-C60-C61-C62
26	T	104	CDL	C51-C52-C53-C54
27	C	306	PEK	C23-C24-C25-C26
20	C	307	PGV	C2-C3-C4-C5
26	G	102	CDL	C71-C72-C73-C74
19	Y	101	TGL	OA1-CA1-OG1-CG1
19	D	201	TGL	C10-C11-C12-C13
19	L	101	TGL	C18-C19-C33-C34
19	A	607	TGL	CA6-CA7-CA8-CA9
19	Y	101	TGL	CB4-CB5-CB6-CB7
19	Y	101	TGL	C19-C33-C34-C35
26	G	102	CDL	C71-CB7-OB8-CB6
26	G	102	CDL	CA3-CA4-CA6-OA8
19	A	607	TGL	C22-C23-C24-C25
19	O	302	TGL	CB2-CB3-CB4-CB5
19	Y	101	TGL	CA5-CA6-CA7-CA8
26	P	304	CDL	C31-C32-C33-C34
27	T	101	PEK	C1-C2-C3-C4
19	L	101	TGL	C11-C12-C13-C14
19	Q	201	TGL	C16-C15-CC9-CC8
19	Y	101	TGL	CB2-CB3-CB4-CB5
19	Y	101	TGL	C10-C11-C12-C13
20	N	607	PGV	C22-C23-C24-C25
20	N	608	PGV	C28-C29-C30-C31
20	N	608	PGV	C30-C31-C32-C33
24	B	303	PSC	C3-C4-C5-C6
24	O	303	PSC	C21-C22-C23-C24
27	C	306	PEK	C16-C17-C18-C19
19	A	607	TGL	CC4-CC5-CC6-CC7
19	L	101	TGL	C22-C23-C24-C25
26	P	304	CDL	C73-C74-C75-C76
21	A	612	EDO	O1-C1-C2-O2
21	A	613	EDO	O1-C1-C2-O2
21	B	304	EDO	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
19	Y	101	TGL	C11-C10-CB9-CB8
19	A	607	TGL	CB2-CB3-CB4-CB5
19	Y	101	TGL	C13-C14-C29-C30
24	O	303	PSC	C23-C24-C25-C26
27	G	101	PEK	C30-C31-C32-C33
26	G	102	CDL	CA7-C31-C32-C33
19	O	302	TGL	CA4-CA5-CA6-CA7
20	N	608	PGV	C26-C27-C28-C29
26	P	304	CDL	C83-C84-C85-C86
26	G	102	CDL	C1-CB2-OB2-PB2
27	G	101	PEK	C10-C11-C12-C13
19	A	607	TGL	CA5-CA6-CA7-CA8
19	Y	101	TGL	CC5-CC6-CC7-CC8
19	A	607	TGL	CB9-C10-C11-C12
19	Q	201	TGL	CC2-CC3-CC4-CC5
20	P	303	PGV	C14-C15-C16-C17
26	G	102	CDL	C51-C52-C53-C54
19	A	607	TGL	C12-C13-C14-C29
20	P	303	PGV	C24-C25-C26-C27
26	C	303	CDL	C52-C53-C54-C55
27	C	306	PEK	O04-C21-O03-C01
20	C	302	PGV	C7-C8-C9-C10
20	P	303	PGV	C25-C26-C27-C28
26	T	104	CDL	C33-C34-C35-C36
27	G	103	PEK	C25-C26-C27-C28
27	C	306	PEK	O02-C1-O01-C02
20	C	307	PGV	C30-C31-C32-C33
19	L	101	TGL	CC3-CC4-CC5-CC6
19	Q	201	TGL	CA4-CA5-CA6-CA7
19	Y	101	TGL	CB7-CB8-CB9-C10
20	C	307	PGV	C3-C4-C5-C6
20	C	307	PGV	C23-C24-C25-C26
19	O	302	TGL	CA6-CA7-CA8-CA9
24	O	303	PSC	C4-C5-C6-C7
26	G	102	CDL	C60-C61-C62-C63
26	P	304	CDL	C55-C56-C57-C58
19	D	201	TGL	C16-C15-CC9-CC8
19	O	302	TGL	C13-C14-C29-C30
20	A	609	PGV	C29-C30-C31-C32
26	T	104	CDL	C37-C38-C39-C40
19	Y	101	TGL	CA6-CA7-CA8-CA9
20	C	307	PGV	C20-C21-C22-C23

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Mol	Chain	Res	Type	Atoms
26	G	102	CDL	C75-C76-C77-C78
19	L	101	TGL	C16-C15-CC9-CC8
19	O	302	TGL	CB9-C10-C11-C12
19	Q	201	TGL	CA2-CA3-CA4-CA5
26	C	303	CDL	C42-C43-C44-C45
27	C	306	PEK	C25-C26-C27-C28
19	A	607	TGL	C11-C12-C13-C14
19	D	201	TGL	C16-C17-C18-C19
19	O	302	TGL	CA2-CA3-CA4-CA5
20	A	609	PGV	C4-C5-C6-C7
24	O	303	PSC	C3-C4-C5-C6
26	C	303	CDL	C36-C37-C38-C39
26	T	104	CDL	C39-C40-C41-C42
20	A	609	PGV	C24-C25-C26-C27
26	C	303	CDL	C60-C61-C62-C63
24	O	303	PSC	C2-C1-O01-C02
27	C	306	PEK	C2-C1-O01-C02
20	C	307	PGV	C1-C2-C3-C4
19	O	302	TGL	CB4-CB5-CB6-CB7
27	T	101	PEK	C34-C35-C36-C37
20	A	609	PGV	C12-C13-C14-C15
20	N	608	PGV	C11-C10-C9-C8
24	O	303	PSC	C13-C14-C15-C16
27	T	103	PEK	C2-C3-C4-C5
27	T	103	PEK	C15-C16-C17-C18
19	D	201	TGL	C17-C18-C19-C33
20	P	301	PGV	C22-C23-C24-C25
26	T	104	CDL	C59-C60-C61-C62
20	P	301	PGV	C6-C7-C8-C9
19	O	302	TGL	CA3-CA4-CA5-CA6
19	Q	201	TGL	C13-C14-C29-C30
20	N	607	PGV	C20-C21-C22-C23
27	T	102	PEK	C30-C31-C32-C33
24	B	303	PSC	C27-C28-C29-C30
27	G	101	PEK	C24-C25-C26-C27
20	C	302	PGV	C10-C11-C12-C13
20	C	302	PGV	C23-C24-C25-C26
29	M	101	DMU	C25-C28-C31-C34
23	J	101	CHD	C13-C17-C20-C21
26	G	102	CDL	C83-C84-C85-C86
23	C	305	CHD	C16-C17-C20-C22
26	G	102	CDL	C57-C58-C59-C60

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Mol	Chain	Res	Type	Atoms
19	O	302	TGL	OB1-CB1-OG2-CG2
19	L	101	TGL	C12-C13-C14-C29
27	T	103	PEK	C23-C24-C25-C26
19	A	607	TGL	CB6-CB7-CB8-CB9
19	L	101	TGL	CA7-CA8-CA9-C20
19	L	101	TGL	CB3-CB4-CB5-CB6
26	P	304	CDL	C16-C17-C18-C19
26	P	304	CDL	C63-C64-C65-C66
19	O	302	TGL	C12-C13-C14-C29
26	T	104	CDL	C35-C36-C37-C38
27	T	103	PEK	C27-C28-C29-C30
20	P	301	PGV	O03-C01-C02-O01
24	B	303	PSC	C13-C14-C15-C16
27	G	101	PEK	C15-C16-C17-C18
27	T	102	PEK	C2-C3-C4-C5
19	L	101	TGL	CC6-CC7-CC8-CC9
19	Y	101	TGL	CA4-CA5-CA6-CA7
19	Y	101	TGL	CC7-CC8-CC9-C15
19	A	607	TGL	CA2-CA3-CA4-CA5
19	Y	101	TGL	CC4-CC5-CC6-CC7
26	G	102	CDL	C12-C13-C14-C15
19	D	201	TGL	CB2-CB3-CB4-CB5
19	O	302	TGL	CB7-CB8-CB9-C10
26	G	102	CDL	C74-C75-C76-C77
20	A	608	PGV	C22-C23-C24-C25
20	N	607	PGV	C21-C22-C23-C24
26	G	102	CDL	C21-C22-C23-C24
27	T	101	PEK	C23-C24-C25-C26
24	B	303	PSC	C02-C03-O11-P
24	O	303	PSC	C5-C6-C7-C8
26	G	102	CDL	C63-C64-C65-C66
26	T	104	CDL	C58-C59-C60-C61
20	C	302	PGV	C14-C15-C16-C17
26	C	303	CDL	C40-C41-C42-C43
26	G	102	CDL	C58-C59-C60-C61
27	G	103	PEK	C22-C23-C24-C25
19	L	101	TGL	CB2-CB3-CB4-CB5
26	G	102	CDL	C72-C73-C74-C75
19	A	607	TGL	CC3-CC4-CC5-CC6
27	T	102	PEK	C4-C5-C6-C7
19	A	607	TGL	CC5-CC6-CC7-CC8
26	P	304	CDL	C82-C83-C84-C85

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Mol	Chain	Res	Type	Atoms
27	C	306	PEK	C33-C34-C35-C36
26	G	102	CDL	OB9-CB7-OB8-CB6
19	A	607	TGL	CA3-CA4-CA5-CA6
23	J	101	CHD	C16-C17-C20-C21
27	G	103	PEK	C33-C34-C35-C36
23	W	101	CHD	C13-C17-C20-C21
20	P	301	PGV	C30-C31-C32-C33
24	B	303	PSC	C23-C24-C25-C26
26	T	104	CDL	C62-C63-C64-C65
19	Y	101	TGL	C20-C21-C22-C23
26	P	304	CDL	C13-C14-C15-C16
20	A	608	PGV	C12-C13-C14-C15
19	O	302	TGL	CC5-CC6-CC7-CC8
20	A	609	PGV	C2-C3-C4-C5
20	C	307	PGV	C27-C28-C29-C30
20	A	609	PGV	C01-C02-C03-O11
26	C	303	CDL	OA5-CA3-CA4-CA6
27	T	103	PEK	C01-C02-C03-O11
24	O	303	PSC	O02-C1-O01-C02
19	Y	101	TGL	CC9-C15-C16-C17
26	G	102	CDL	C52-C53-C54-C55
19	A	607	TGL	C18-C19-C33-C34
24	B	303	PSC	C24-C25-C26-C27
26	G	102	CDL	C11-C12-C13-C14
26	C	303	CDL	C61-C62-C63-C64
26	C	303	CDL	C71-C72-C73-C74
27	G	101	PEK	C34-C35-C36-C37
27	T	103	PEK	C16-C17-C18-C19
19	A	607	TGL	CA1-CA2-CA3-CA4
19	O	302	TGL	CA1-CA2-CA3-CA4
20	A	609	PGV	C19-C20-C21-C22
20	A	609	PGV	C13-C14-C15-C16
19	A	607	TGL	C11-C10-CB9-CB8
19	O	302	TGL	C21-C20-CA9-CA8
26	G	102	CDL	C13-C14-C15-C16
26	P	304	CDL	C41-C42-C43-C44
26	P	304	CDL	C56-C57-C58-C59
20	N	608	PGV	C29-C30-C31-C32
27	C	306	PEK	C26-C27-C28-C29
19	Q	201	TGL	C21-C20-CA9-CA8
19	D	201	TGL	CC5-CC6-CC7-CC8
29	Z	101	DMU	C3-C4-C57-O61

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Mol	Chain	Res	Type	Atoms
19	Q	201	TGL	C23-C24-C25-C26
19	D	201	TGL	OG1-CG1-CG2-CG3
19	Y	101	TGL	OG1-CG1-CG2-CG3
20	C	307	PGV	O03-C01-C02-C03
20	P	301	PGV	O03-C01-C02-C03
26	P	304	CDL	CA3-CA4-CA6-OA8
27	C	306	PEK	O03-C01-C02-C03
27	C	306	PEK	C15-C16-C17-C18
19	Q	201	TGL	C19-C33-C34-C35
26	C	303	CDL	C76-C77-C78-C79
26	T	104	CDL	C63-C64-C65-C66
19	A	607	TGL	C19-C33-C34-C35
19	D	201	TGL	CC6-CC7-CC8-CC9
20	A	609	PGV	C26-C27-C28-C29
20	P	303	PGV	C27-C28-C29-C30
21	T	105	EDO	O1-C1-C2-O2
19	L	101	TGL	CA4-CA5-CA6-CA7
27	T	103	PEK	C30-C31-C32-C33
19	Q	201	TGL	CA9-C20-C21-C22
26	P	304	CDL	C61-C62-C63-C64
27	T	102	PEK	C32-C33-C34-C35
26	P	304	CDL	C52-C53-C54-C55
19	O	302	TGL	CB6-CB7-CB8-CB9
26	C	303	CDL	C77-C78-C79-C80
26	C	303	CDL	C23-C24-C25-C26
26	G	102	CDL	C23-C24-C25-C26
26	P	304	CDL	C79-C80-C81-C82
27	G	101	PEK	C23-C24-C25-C26
19	Q	201	TGL	CA3-CA4-CA5-CA6
19	Y	101	TGL	C16-C15-CC9-CC8
26	G	102	CDL	C35-C36-C37-C38
19	D	201	TGL	CA2-CA3-CA4-CA5
26	C	303	CDL	C75-C76-C77-C78
20	A	609	PGV	C03-C02-O01-C1
24	O	303	PSC	C01-C02-O01-C1
19	D	201	TGL	C20-C21-C22-C23
19	D	201	TGL	C21-C22-C23-C24
19	Q	201	TGL	CB7-CB8-CB9-C10
24	O	303	PSC	C28-C29-C30-C31
26	C	303	CDL	C17-C18-C19-C20
20	A	608	PGV	C11-C10-C9-C8
20	C	307	PGV	C12-C13-C14-C15

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Mol	Chain	Res	Type	Atoms
20	P	301	PGV	C12-C13-C14-C15
20	P	303	PGV	C12-C13-C14-C15
19	O	302	TGL	C16-C17-C18-C19
20	A	609	PGV	C25-C26-C27-C28
27	T	103	PEK	C25-C26-C27-C28
14	N	602	HEA	C2A-C3A-CMA-OMA
26	C	303	CDL	C11-C12-C13-C14
27	G	103	PEK	C31-C32-C33-C34
27	T	101	PEK	C4-C5-C6-C7
26	C	303	CDL	C37-C38-C39-C40
27	G	103	PEK	O01-C02-C03-O11
19	A	607	TGL	C13-C14-C29-C30
20	A	609	PGV	C7-C8-C9-C10
20	P	303	PGV	C29-C30-C31-C32
26	P	304	CDL	C35-C36-C37-C38
19	A	607	TGL	C23-C24-C25-C26
20	C	307	PGV	C4-C5-C6-C7
27	T	102	PEK	C35-C36-C37-C38
20	A	608	PGV	C14-C15-C16-C17
20	N	607	PGV	C6-C7-C8-C9
26	P	304	CDL	C37-C38-C39-C40
29	Z	101	DMU	O16-C18-C19-C22
19	A	607	TGL	C21-C20-CA9-CA8
20	C	302	PGV	C22-C23-C24-C25
26	C	303	CDL	C74-C75-C76-C77
19	A	607	TGL	CB4-CB5-CB6-CB7
19	D	201	TGL	CA7-CA8-CA9-C20
20	C	307	PGV	C21-C22-C23-C24
26	G	102	CDL	C16-C17-C18-C19
27	T	102	PEK	C25-C26-C27-C28
19	Q	201	TGL	C20-C21-C22-C23
27	G	101	PEK	C25-C26-C27-C28
20	C	307	PGV	O03-C01-C02-O01
26	G	102	CDL	C19-C20-C21-C22
29	M	101	DMU	C28-C31-C34-C37
26	C	303	CDL	C84-C85-C86-C87
19	Q	201	TGL	CC4-CC5-CC6-CC7
20	N	607	PGV	C27-C28-C29-C30
26	T	104	CDL	C72-C73-C74-C75
26	C	303	CDL	C62-C63-C64-C65
26	T	104	CDL	C55-C56-C57-C58
26	G	102	CDL	C39-C40-C41-C42

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Mol	Chain	Res	Type	Atoms
20	A	608	PGV	C31-C32-C33-C34
26	P	304	CDL	C24-C25-C26-C27
26	P	304	CDL	C84-C85-C86-C87
26	T	104	CDL	C24-C25-C26-C27
19	O	302	TGL	C25-C26-C27-C28
27	T	103	PEK	C17-C18-C19-C20
19	Q	201	TGL	CC6-CC7-CC8-CC9
19	D	201	TGL	CB4-CB5-CB6-CB7
26	C	303	CDL	C33-C34-C35-C36
19	Q	201	TGL	CA1-CA2-CA3-CA4
19	A	607	TGL	C29-C30-C31-C32
24	O	303	PSC	C11-C10-C9-C8
20	P	301	PGV	C31-C32-C33-C34
26	G	102	CDL	C44-C45-C46-C47
26	T	104	CDL	C40-C41-C42-C43
14	N	602	HEA	C4D-C3D-CAD-CBD
26	P	304	CDL	C77-C78-C79-C80
23	G	104	CHD	C17-C20-C22-C23
26	C	303	CDL	C18-C19-C20-C21
20	C	307	PGV	C02-C03-O11-P
20	A	608	PGV	C23-C24-C25-C26
19	O	302	TGL	C11-C12-C13-C14
26	T	104	CDL	C81-C82-C83-C84
24	O	303	PSC	C31-C32-C33-C34
26	P	304	CDL	OA5-CA3-CA4-CA6
24	O	303	PSC	C22-C23-C24-C25
19	O	302	TGL	C20-C21-C22-C23
27	T	101	PEK	C25-C26-C27-C28
19	A	607	TGL	CB5-CB6-CB7-CB8
19	L	101	TGL	C21-C20-CA9-CA8
26	G	102	CDL	C17-C18-C19-C20
26	P	304	CDL	C80-C81-C82-C83
26	C	303	CDL	C35-C36-C37-C38
19	O	302	TGL	C29-C30-C31-C32
19	Q	201	TGL	CA5-CA6-CA7-CA8
19	L	101	TGL	OG1-CG1-CG2-CG3
26	C	303	CDL	CB3-CB4-CB6-OB8
26	P	304	CDL	CB3-CB4-CB6-OB8
26	T	104	CDL	CA3-CA4-CA6-OA8
19	A	607	TGL	C25-C26-C27-C28
20	A	609	PGV	C22-C23-C24-C25
20	N	607	PGV	C29-C30-C31-C32

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Mol	Chain	Res	Type	Atoms
20	P	301	PGV	C13-C14-C15-C16
26	G	102	CDL	C43-C44-C45-C46
19	L	101	TGL	C13-C14-C29-C30
26	T	104	CDL	C14-C15-C16-C17
21	K	102	EDO	O1-C1-C2-O2
21	P	308	EDO	O1-C1-C2-O2
19	O	302	TGL	CC4-CC5-CC6-CC7
19	O	302	TGL	CC6-CC7-CC8-CC9
19	Q	201	TGL	C10-C11-C12-C13
20	P	301	PGV	O01-C02-C03-O11
27	T	102	PEK	O01-C02-C03-O11
20	C	302	PGV	C25-C26-C27-C28
14	N	602	HEA	C2D-C3D-CAD-CBD
26	C	303	CDL	C78-C79-C80-C81
26	P	304	CDL	C14-C15-C16-C17
23	B	302	CHD	C17-C20-C22-C23
20	A	608	PGV	O03-C19-C20-C21
19	Y	101	TGL	OG1-CG1-CG2-OG2
20	A	609	PGV	O03-C01-C02-O01
24	B	303	PSC	O03-C01-C02-O01
24	B	303	PSC	C10-C11-C12-C13
24	O	303	PSC	C9-C10-C11-C12
24	O	303	PSC	C10-C11-C12-C13
27	C	306	PEK	C5-C6-C7-C8
27	C	306	PEK	C6-C7-C8-C9
27	C	306	PEK	C11-C10-C9-C8
27	G	103	PEK	C11-C10-C9-C8
27	G	103	PEK	C9-C10-C11-C12
27	G	103	PEK	C12-C13-C14-C15
27	T	101	PEK	C11-C10-C9-C8
27	T	102	PEK	C9-C10-C11-C12
27	T	103	PEK	C6-C7-C8-C9
27	T	103	PEK	C11-C10-C9-C8
27	T	103	PEK	C12-C13-C14-C15
20	P	301	PGV	C25-C26-C27-C28
20	P	303	PGV	C31-C32-C33-C34
19	D	201	TGL	CA3-CA4-CA5-CA6
19	Q	201	TGL	C24-C25-C26-C27
26	P	304	CDL	C59-C60-C61-C62
19	A	607	TGL	C33-C34-C35-C36
19	D	201	TGL	CB5-CB6-CB7-CB8
26	P	304	CDL	C22-C23-C24-C25

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Mol	Chain	Res	Type	Atoms
27	G	101	PEK	C32-C33-C34-C35
19	Q	201	TGL	C15-C16-C17-C18
27	G	101	PEK	C33-C34-C35-C36
19	Q	201	TGL	CC5-CC6-CC7-CC8
19	O	302	TGL	CC7-CC8-CC9-C15
20	C	307	PGV	C14-C15-C16-C17
26	T	104	CDL	C22-C23-C24-C25
27	T	101	PEK	C17-C18-C19-C20
19	Q	201	TGL	CB3-CB4-CB5-CB6
27	G	101	PEK	C26-C27-C28-C29
19	O	302	TGL	C16-C15-CC9-CC8
19	Y	101	TGL	CC1-CC2-CC3-CC4
26	P	304	CDL	CA5-C11-C12-C13
20	N	608	PGV	C31-C32-C33-C34
26	T	104	CDL	C41-C42-C43-C44
26	G	102	CDL	C15-C16-C17-C18
27	G	103	PEK	C34-C35-C36-C37
29	M	101	DMU	C19-C22-C25-C28
20	N	607	PGV	C01-C02-C03-O11
20	P	301	PGV	C01-C02-C03-O11
24	B	303	PSC	C01-C02-C03-O11
27	G	103	PEK	C01-C02-C03-O11
27	T	102	PEK	C01-C02-C03-O11
19	Y	101	TGL	CB5-CB6-CB7-CB8
26	G	102	CDL	C62-C63-C64-C65
19	L	101	TGL	CA9-C20-C21-C22
20	P	303	PGV	C13-C14-C15-C16
20	P	301	PGV	C3-C4-C5-C6
20	A	609	PGV	C31-C32-C33-C34
19	A	607	TGL	CB1-CB2-CB3-CB4
27	T	101	PEK	C7-C8-C9-C10
27	T	103	PEK	O02-C1-O01-C02
27	G	101	PEK	C16-C17-C18-C19
26	P	304	CDL	C34-C35-C36-C37
19	Y	101	TGL	CC3-CC4-CC5-CC6
19	A	607	TGL	C24-C25-C26-C27
20	N	607	PGV	C12-C13-C14-C15
24	O	303	PSC	C6-C7-C8-C9
27	T	101	PEK	C15-C16-C17-C18
19	L	101	TGL	C17-C18-C19-C33
27	G	103	PEK	C27-C28-C29-C30
20	N	607	PGV	C23-C24-C25-C26

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Mol	Chain	Res	Type	Atoms
20	A	609	PGV	O01-C02-C03-O11
20	N	607	PGV	O01-C02-C03-O11
27	T	103	PEK	O01-C02-C03-O11
19	Q	201	TGL	CB4-CB5-CB6-CB7
19	Y	101	TGL	CG1-CG2-CG3-OG3
24	O	303	PSC	O03-C01-C02-C03
27	T	103	PEK	O03-C01-C02-C03
27	T	103	PEK	C2-C1-O01-C02
26	P	304	CDL	C51-C52-C53-C54
29	M	101	DMU	C22-C25-C28-C31
26	G	102	CDL	CB5-C51-C52-C53
26	T	104	CDL	C71-C72-C73-C74
21	K	101	EDO	O1-C1-C2-O2
21	L	102	EDO	O1-C1-C2-O2
19	D	201	TGL	OG1-CG1-CG2-OG2
19	Y	101	TGL	OG2-CG2-CG3-OG3
24	O	303	PSC	O03-C01-C02-O01
26	C	303	CDL	OB6-CB4-CB6-OB8
27	T	102	PEK	O03-C01-C02-O01
19	D	201	TGL	C13-C14-C29-C30
27	T	103	PEK	C26-C27-C28-C29
27	T	103	PEK	C28-C29-C30-C31
20	A	608	PGV	C30-C31-C32-C33
26	C	303	CDL	C58-C59-C60-C61
20	P	301	PGV	C4-C5-C6-C7
26	T	104	CDL	CB7-C71-C72-C73
27	T	102	PEK	C2-C1-O01-C02
20	P	301	PGV	C02-C03-O11-P
26	P	304	CDL	C1-CB2-OB2-PB2
26	C	303	CDL	CB4-CB6-OB8-CB7
27	C	306	PEK	C34-C35-C36-C37
20	N	608	PGV	C19-C20-C21-C22
19	L	101	TGL	C20-C21-C22-C23
19	D	201	TGL	CC9-C15-C16-C17
19	A	607	TGL	CA4-CA5-CA6-CA7
24	B	303	PSC	C15-C16-C17-C18
26	C	303	CDL	C32-C33-C34-C35
27	C	306	PEK	C30-C31-C32-C33
24	B	303	PSC	C31-C32-C33-C34
27	G	103	PEK	C3-C4-C5-C6
26	P	304	CDL	CA4-CA3-OA5-PA1
26	T	104	CDL	C1-CB2-OB2-PB2

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Mol	Chain	Res	Type	Atoms
20	P	303	PGV	C28-C29-C30-C31
24	B	303	PSC	O01-C02-C03-O11
26	C	303	CDL	OA5-CA3-CA4-OA6
27	T	101	PEK	C27-C28-C29-C30
14	N	601	HEA	C4A-C3A-CMA-OMA
14	N	602	HEA	C4A-C3A-CMA-OMA
19	O	302	TGL	CA5-CA6-CA7-CA8
20	N	608	PGV	C5-C6-C7-C8
26	G	102	CDL	OA6-CA4-CA6-OA8
26	P	304	CDL	OA6-CA4-CA6-OA8
26	P	304	CDL	OB6-CB4-CB6-OB8
26	T	104	CDL	OA6-CA4-CA6-OA8
20	A	609	PGV	O03-C01-C02-C03
20	N	607	PGV	O03-C01-C02-C03
26	C	303	CDL	CA3-CA4-CA6-OA8
19	Y	101	TGL	CA2-CA3-CA4-CA5
19	Y	101	TGL	C17-C18-C19-C33
27	G	103	PEK	C28-C29-C30-C31
24	O	303	PSC	C15-C16-C17-C18
19	L	101	TGL	CA2-CA3-CA4-CA5
20	N	607	PGV	C31-C32-C33-C34
24	B	303	PSC	C2-C3-C4-C5
20	A	609	PGV	C04-O12-P-O13
20	C	307	PGV	C04-O12-P-O13
20	P	301	PGV	C03-O11-P-O13
20	P	301	PGV	C04-O12-P-O11
20	P	303	PGV	C04-O12-P-O14
24	B	303	PSC	C03-O11-P-O14
24	B	303	PSC	C04-O12-P-O11
24	B	303	PSC	C04-O12-P-O13
26	C	303	CDL	CA3-OA5-PA1-OA3
26	G	102	CDL	CB2-OB2-PB2-OB3
26	G	102	CDL	CB2-OB2-PB2-OB4
26	G	102	CDL	CB3-OB5-PB2-OB3
26	P	304	CDL	CB3-OB5-PB2-OB3
26	T	104	CDL	CA2-OA2-PA1-OA5
26	T	104	CDL	CA3-OA5-PA1-OA3
27	T	102	PEK	O12-C04-C05-N
19	Y	101	TGL	C22-C23-C24-C25
26	C	303	CDL	C31-C32-C33-C34
20	C	302	PGV	C02-C03-O11-P
20	P	303	PGV	C02-C03-O11-P

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Mol	Chain	Res	Type	Atoms
26	C	303	CDL	C1-CA2-OA2-PA1
26	C	303	CDL	CA4-CA3-OA5-PA1
27	T	102	PEK	C02-C03-O11-P
19	D	201	TGL	CB7-CB8-CB9-C10
19	D	201	TGL	C14-C29-C30-C31
26	P	304	CDL	C39-C40-C41-C42
26	P	304	CDL	C78-C79-C80-C81
20	C	307	PGV	O05-C05-C06-O06
26	P	304	CDL	C74-C75-C76-C77
26	P	304	CDL	C72-C73-C74-C75
27	G	101	PEK	C21-C22-C23-C24
19	D	201	TGL	CA4-CA5-CA6-CA7
19	L	101	TGL	C10-C11-C12-C13
26	G	102	CDL	C64-C65-C66-C67
27	C	306	PEK	C28-C29-C30-C31
20	A	608	PGV	C11-C12-C13-C14
26	G	102	CDL	CB7-C71-C72-C73
20	N	608	PGV	C27-C28-C29-C30
20	N	608	PGV	O03-C19-C20-C21
20	A	608	PGV	C24-C25-C26-C27
20	P	303	PGV	C4-C5-C6-C7
20	P	301	PGV	C23-C24-C25-C26
27	T	102	PEK	O02-C1-O01-C02
26	P	304	CDL	C57-C58-C59-C60
19	L	101	TGL	OG1-CG1-CG2-OG2
19	D	201	TGL	CA9-C20-C21-C22
24	O	303	PSC	O03-C19-C20-C21
27	T	101	PEK	C35-C36-C37-C38
24	B	303	PSC	O03-C01-C02-C03
26	G	102	CDL	C81-C82-C83-C84
20	P	301	PGV	C27-C28-C29-C30
20	A	608	PGV	C21-C22-C23-C24
21	O	304	EDO	O1-C1-C2-O2
19	L	101	TGL	C29-C30-C31-C32
27	T	102	PEK	C31-C32-C33-C34
27	T	103	PEK	C29-C30-C31-C32
19	L	101	TGL	CA5-CA6-CA7-CA8
20	C	307	PGV	C11-C10-C9-C8
26	P	304	CDL	C12-C13-C14-C15
19	D	201	TGL	CB9-C10-C11-C12
27	T	101	PEK	C32-C33-C34-C35
19	D	201	TGL	C19-C33-C34-C35

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Mol	Chain	Res	Type	Atoms
26	P	304	CDL	C40-C41-C42-C43
26	T	104	CDL	C75-C76-C77-C78
26	P	304	CDL	OB9-CB7-OB8-CB6
26	T	104	CDL	C77-C78-C79-C80
20	N	607	PGV	C11-C12-C13-C14
20	N	608	PGV	C9-C10-C11-C12
19	A	607	TGL	C14-C29-C30-C31
26	G	102	CDL	OB6-CB4-CB6-OB8
26	G	102	CDL	C54-C55-C56-C57
26	G	102	CDL	C79-C80-C81-C82
19	Y	101	TGL	CA7-CA8-CA9-C20
26	C	303	CDL	C63-C64-C65-C66
26	P	304	CDL	C43-C44-C45-C46
26	P	304	CDL	C71-CB7-OB8-CB6
19	A	607	TGL	OG1-CG1-CG2-CG3
20	P	303	PGV	C15-C16-C17-C18
27	T	102	PEK	C27-C28-C29-C30
19	O	302	TGL	C22-C23-C24-C25
27	T	102	PEK	C24-C25-C26-C27
24	B	303	PSC	C20-C21-C22-C23
20	A	608	PGV	C29-C30-C31-C32
21	B	305	EDO	O1-C1-C2-O2
21	C	308	EDO	O1-C1-C2-O2
19	O	302	TGL	C15-C16-C17-C18
20	A	609	PGV	O03-C19-C20-C21
23	B	302	CHD	C22-C23-C24-O26
23	G	104	CHD	C22-C23-C24-O26
23	J	101	CHD	C16-C17-C20-C22
27	T	102	PEK	C33-C34-C35-C36
19	A	607	TGL	CC7-CC8-CC9-C15
14	A	601	HEA	CAD-CBD-CGD-O1D
14	N	601	HEA	CAD-CBD-CGD-O2D
14	N	602	HEA	CAD-CBD-CGD-O1D
23	B	302	CHD	C22-C23-C24-O25
23	G	104	CHD	C22-C23-C24-O25
27	T	101	PEK	C3-C4-C5-C6
24	O	303	PSC	C02-C03-O11-P
26	T	104	CDL	C31-C32-C33-C34
23	J	101	CHD	C22-C23-C24-O25
20	P	303	PGV	C22-C23-C24-C25
27	C	306	PEK	C1-C2-C3-C4
27	T	103	PEK	C14-C15-C16-C17

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Mol	Chain	Res	Type	Atoms
23	C	304	CHD	C22-C23-C24-O25
24	B	303	PSC	C9-C10-C11-C12
27	G	101	PEK	C6-C7-C8-C9
27	T	103	PEK	C11-C12-C13-C14
27	T	101	PEK	C26-C27-C28-C29
27	C	306	PEK	C31-C32-C33-C34
27	T	102	PEK	C34-C35-C36-C37
26	T	104	CDL	C44-C45-C46-C47
14	N	602	HEA	CAD-CBD-CGD-O2D
19	A	607	TGL	C17-C18-C19-C33
19	D	201	TGL	CC7-CC8-CC9-C15
20	A	609	PGV	C3-C4-C5-C6
20	C	307	PGV	C15-C16-C17-C18
14	N	601	HEA	CAD-CBD-CGD-O1D
20	C	302	PGV	C05-C04-O12-P
20	P	303	PGV	C05-C04-O12-P
19	L	101	TGL	OG1-CA1-CA2-CA3
20	N	608	PGV	C13-C14-C15-C16
27	T	101	PEK	C30-C31-C32-C33
19	L	101	TGL	C14-C29-C30-C31
19	L	101	TGL	CG1-CG2-CG3-OG3
27	T	101	PEK	O03-C01-C02-C03
23	C	304	CHD	C22-C23-C24-O26
19	D	201	TGL	CB6-CB7-CB8-CB9
19	Q	201	TGL	CB5-CB6-CB7-CB8
26	G	102	CDL	OA5-CA3-CA4-OA6
14	N	602	HEA	C26-C15-C16-C17
19	D	201	TGL	C21-C20-CA9-CA8
26	P	304	CDL	C75-C76-C77-C78
21	A	611	EDO	O1-C1-C2-O2
19	L	101	TGL	CC5-CC6-CC7-CC8
26	T	104	CDL	C52-C51-CB5-OB6
26	T	104	CDL	C15-C16-C17-C18
19	Y	101	TGL	C21-C20-CA9-CA8
20	N	607	PGV	C11-C10-C9-C8
27	T	101	PEK	O03-C01-C02-O01
14	A	602	HEA	CAD-CBD-CGD-O1D
26	T	104	CDL	C18-C19-C20-C21
23	J	101	CHD	C22-C23-C24-O26
19	A	607	TGL	CA7-CA8-CA9-C20
23	C	305	CHD	C13-C17-C20-C22
14	A	602	HEA	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
29	Z	101	DMU	O5-C4-C57-O61
26	C	303	CDL	C39-C40-C41-C42
27	T	103	PEK	C31-C32-C33-C34
23	W	101	CHD	C22-C23-C24-O26
19	Q	201	TGL	C18-C19-C33-C34
24	B	303	PSC	C01-C02-O01-C1
27	T	101	PEK	O04-C21-O03-C01
24	O	303	PSC	C7-C8-C9-C10
26	C	303	CDL	C22-C23-C24-C25
26	T	104	CDL	C21-C22-C23-C24
20	N	607	PGV	C24-C25-C26-C27
19	O	302	TGL	C24-C25-C26-C27
19	L	101	TGL	CA3-CA4-CA5-CA6
26	T	104	CDL	OB6-CB4-CB6-OB8
20	P	303	PGV	C9-C10-C11-C12
20	N	608	PGV	C22-C23-C24-C25
19	O	302	TGL	OG3-CC1-CC2-CC3
20	A	609	PGV	O01-C1-C2-C3
20	N	607	PGV	O03-C19-C20-C21
20	P	301	PGV	O01-C1-C2-C3
14	N	602	HEA	C14-C15-C16-C17
27	T	103	PEK	C4-C5-C6-C7
27	T	102	PEK	O01-C1-C2-C3
19	Q	201	TGL	CB9-C10-C11-C12
19	L	101	TGL	C24-C25-C26-C27
19	L	101	TGL	CC4-CC5-CC6-CC7
20	C	302	PGV	C27-C28-C29-C30
24	O	303	PSC	C12-C13-C14-C15
27	T	103	PEK	C3-C4-C5-C6
14	A	601	HEA	CAD-CBD-CGD-O2D
24	B	303	PSC	O03-C19-C20-C21
24	B	303	PSC	C30-C31-C32-C33
19	Q	201	TGL	OG1-CA1-CA2-CA3
26	P	304	CDL	CB5-C51-C52-C53
19	L	101	TGL	C25-C26-C27-C28
26	P	304	CDL	OA5-CA3-CA4-OA6
26	G	102	CDL	C12-C11-CA5-OA6
23	W	101	CHD	C22-C23-C24-O25
19	A	607	TGL	OG3-CC1-CC2-CC3
27	T	101	PEK	O01-C1-C2-C3
19	D	201	TGL	OG1-CA1-CA2-CA3
26	T	104	CDL	C72-C71-CB7-OB8

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Mol	Chain	Res	Type	Atoms
20	A	609	PGV	C21-C22-C23-C24
26	T	104	CDL	C73-C74-C75-C76
26	T	104	CDL	C74-C75-C76-C77
27	G	103	PEK	C15-C16-C17-C18
26	T	104	CDL	C60-C61-C62-C63
20	P	303	PGV	C30-C31-C32-C33
20	A	609	PGV	C05-C04-O12-P
23	P	306	CHD	C22-C23-C24-O26
20	P	301	PGV	C11-C12-C13-C14
19	A	607	TGL	OC1-CC1-CC2-CC3
26	C	303	CDL	CB7-C71-C72-C73
19	Q	201	TGL	OA1-CA1-CA2-CA3
26	G	102	CDL	C12-C11-CA5-OA7
24	B	303	PSC	C22-C23-C24-C25
26	P	304	CDL	C17-C18-C19-C20
19	O	302	TGL	OC1-CC1-CC2-CC3
20	P	301	PGV	O02-C1-C2-C3
24	B	303	PSC	O04-C19-C20-C21
26	C	303	CDL	C43-C44-C45-C46
20	C	307	PGV	C7-C8-C9-C10
19	L	101	TGL	OG2-CG2-CG3-OG3
19	D	201	TGL	OA1-CA1-CA2-CA3
20	N	607	PGV	O04-C19-C20-C21
27	C	306	PEK	O01-C1-C2-C3
27	C	306	PEK	C13-C14-C15-C16
27	T	101	PEK	C22-C21-O03-C01
19	Q	201	TGL	OG1-CG1-CG2-CG3
27	T	102	PEK	O02-C1-C2-C3
26	C	303	CDL	C24-C25-C26-C27
26	T	104	CDL	C72-C71-CB7-OB9
27	C	306	PEK	C24-C25-C26-C27
27	T	101	PEK	O02-C1-C2-C3
20	A	609	PGV	O02-C1-C2-C3
27	T	102	PEK	C17-C18-C19-C20
26	C	303	CDL	C52-C51-CB5-OB6
26	T	104	CDL	C12-C11-CA5-OA6

There are no ring outliers.

48 monomers are involved in 200 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
14	N	602	HEA	5	0

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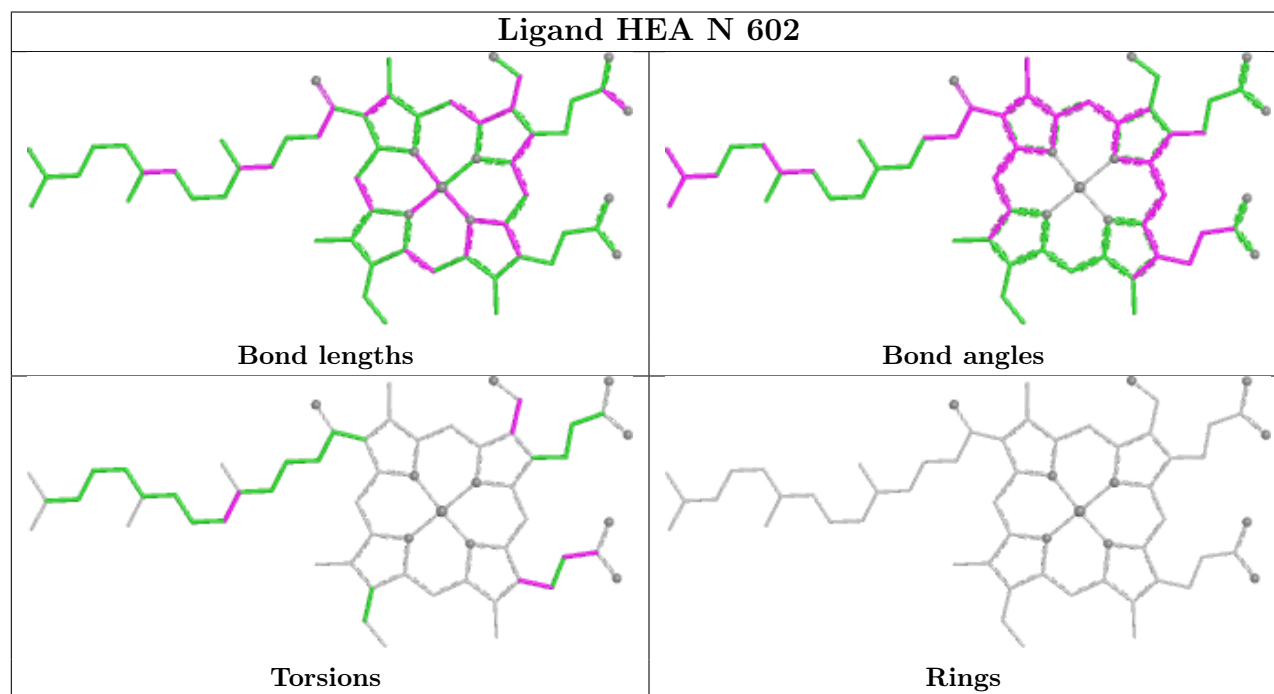
Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	A	606	CMO	1	0
18	N	606	CMO	2	0
27	T	101	PEK	3	0
14	A	601	HEA	6	0
19	L	101	TGL	9	0
20	N	607	PGV	5	0
23	W	101	CHD	9	0
27	T	102	PEK	8	0
19	O	302	TGL	7	0
27	C	306	PEK	3	0
21	C	311	EDO	2	0
20	C	307	PGV	4	0
26	C	303	CDL	6	0
19	Y	101	TGL	5	0
20	P	303	PGV	1	0
21	N	610	EDO	3	0
14	A	602	HEA	8	0
21	F	103	EDO	1	0
27	T	103	PEK	2	0
24	B	303	PSC	4	0
21	T	105	EDO	3	0
20	N	608	PGV	1	0
26	P	304	CDL	7	0
19	D	201	TGL	5	0
21	F	102	EDO	1	0
26	G	102	CDL	6	0
23	G	104	CHD	1	0
21	A	613	EDO	6	0
27	G	103	PEK	3	0
19	Q	201	TGL	5	0
20	A	608	PGV	1	0
21	B	304	EDO	1	0
26	T	104	CDL	21	0
14	N	601	HEA	7	0
23	C	305	CHD	1	0
23	P	306	CHD	3	0
20	P	301	PGV	2	0
29	Z	101	DMU	2	0
19	A	607	TGL	3	0
21	A	612	EDO	12	0
23	P	305	CHD	3	0
23	C	304	CHD	4	0

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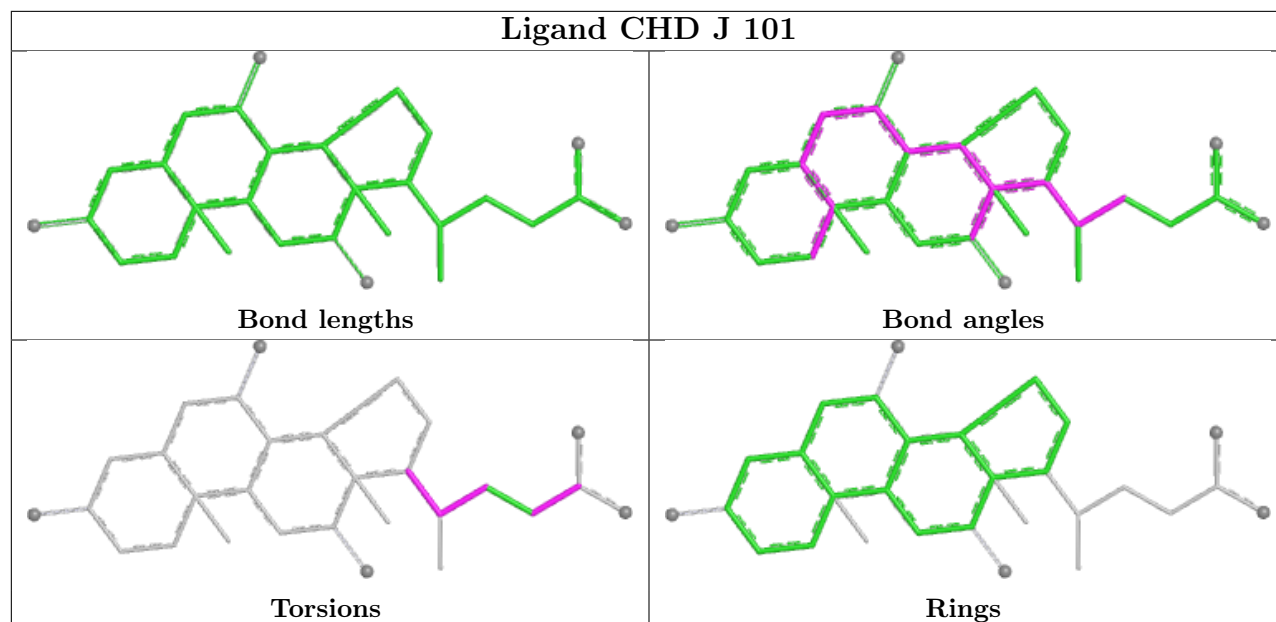
Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
23	B	302	CHD	1	0
24	O	303	PSC	5	0
21	H	101	EDO	2	0
27	G	101	PEK	3	0
20	A	609	PGV	4	0

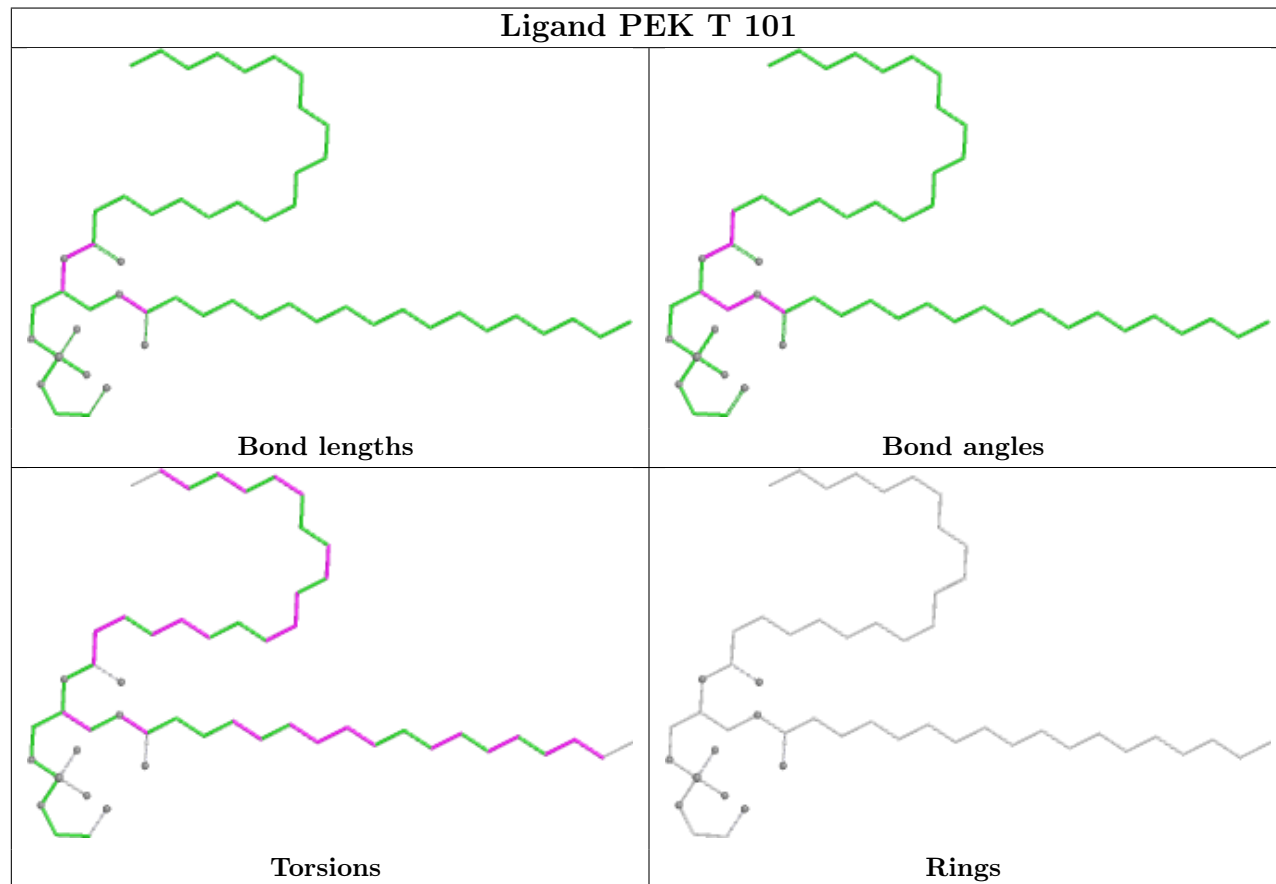
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

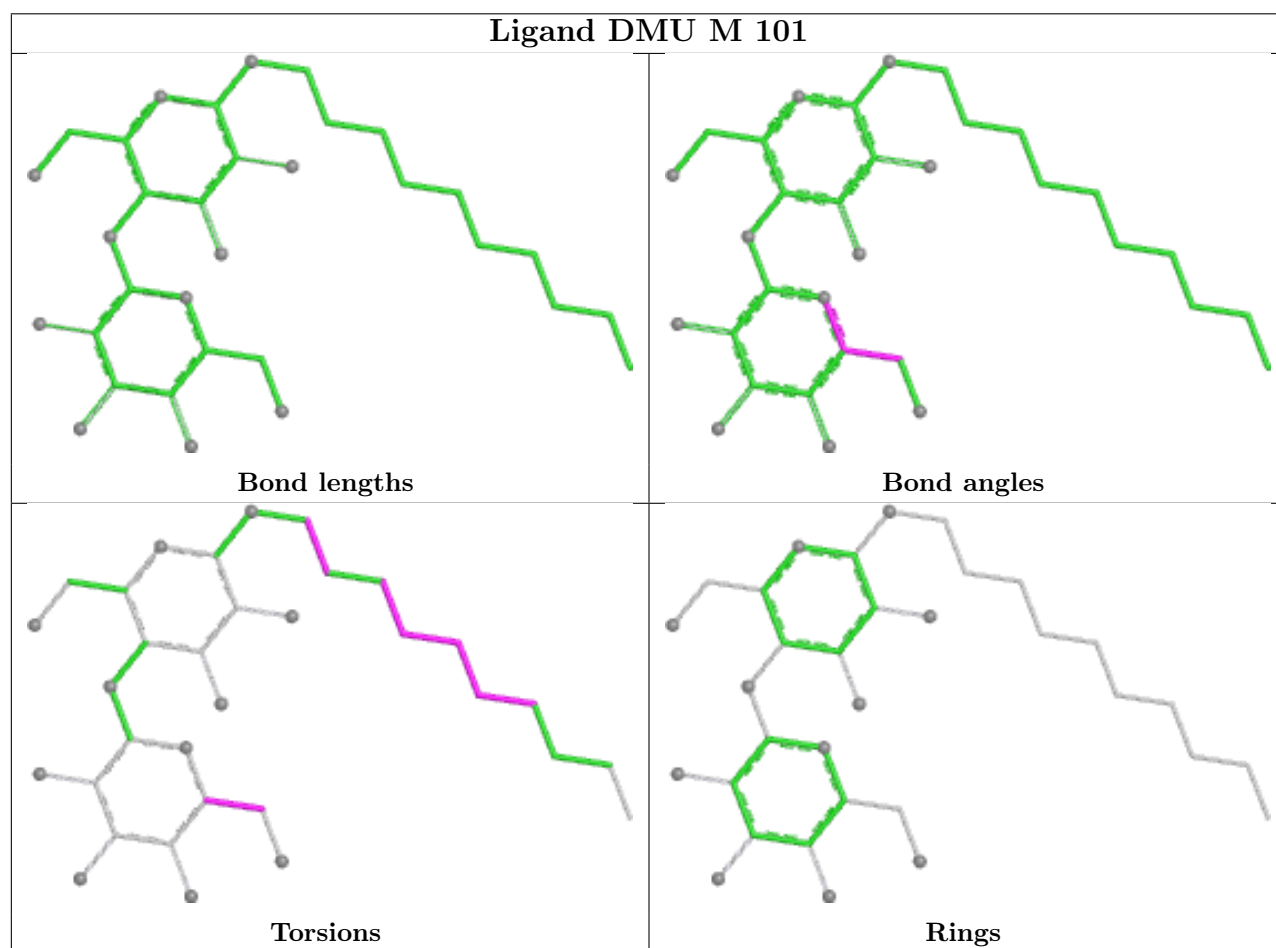
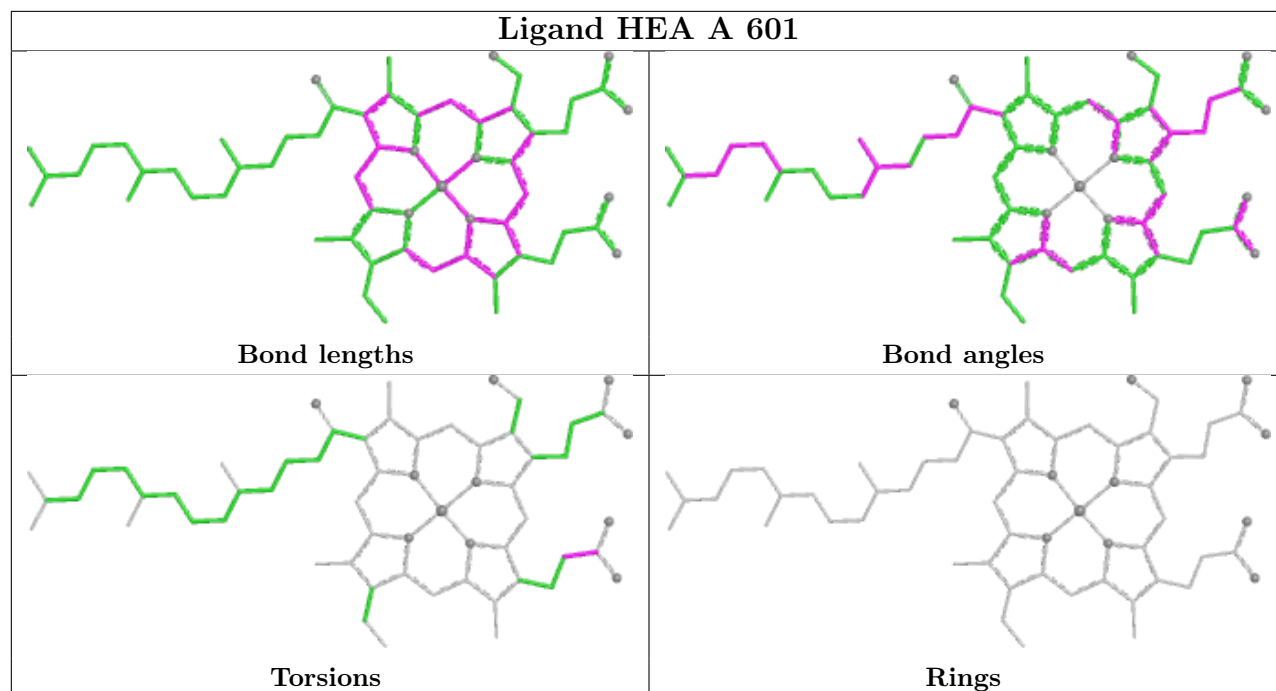


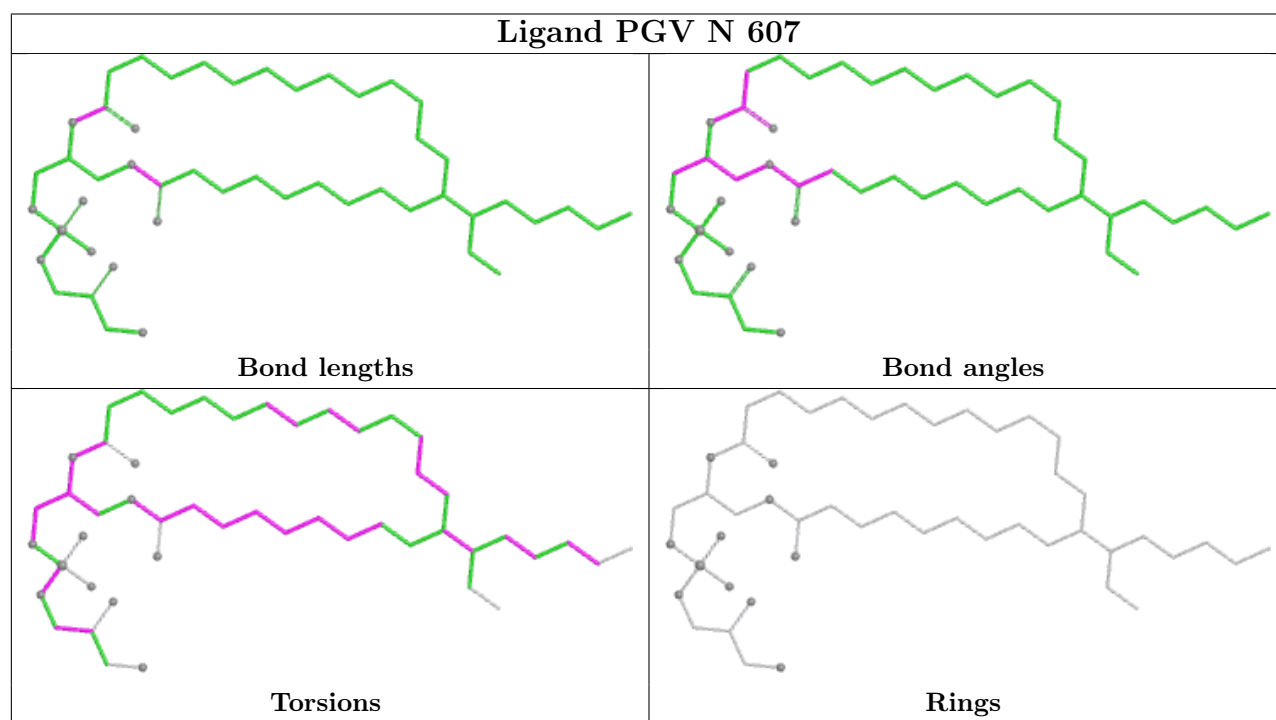
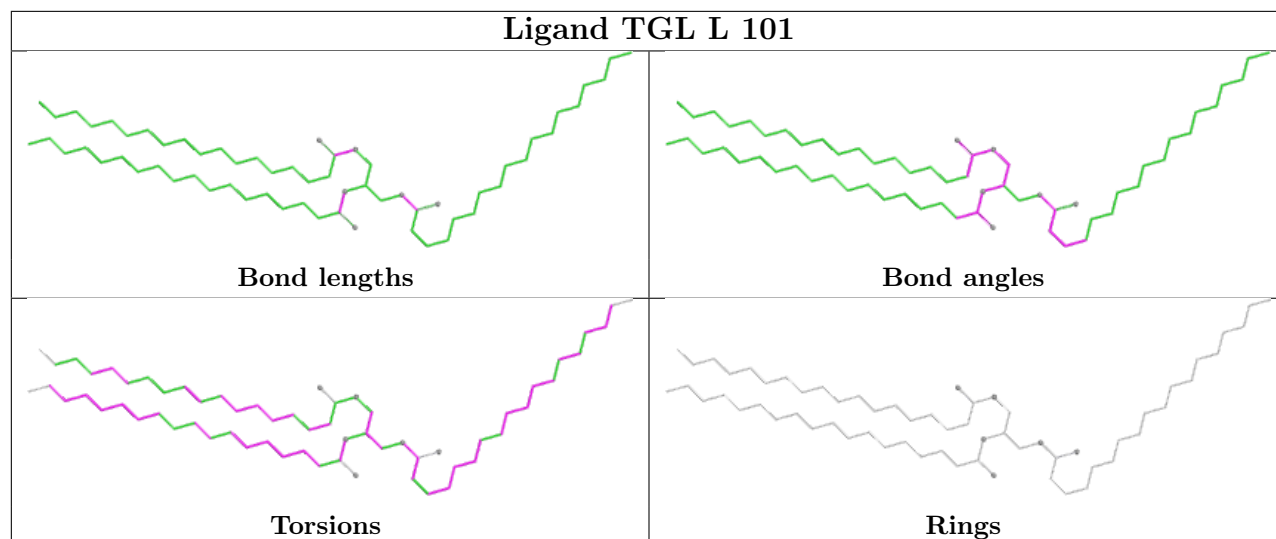
Ligand CHD J 101



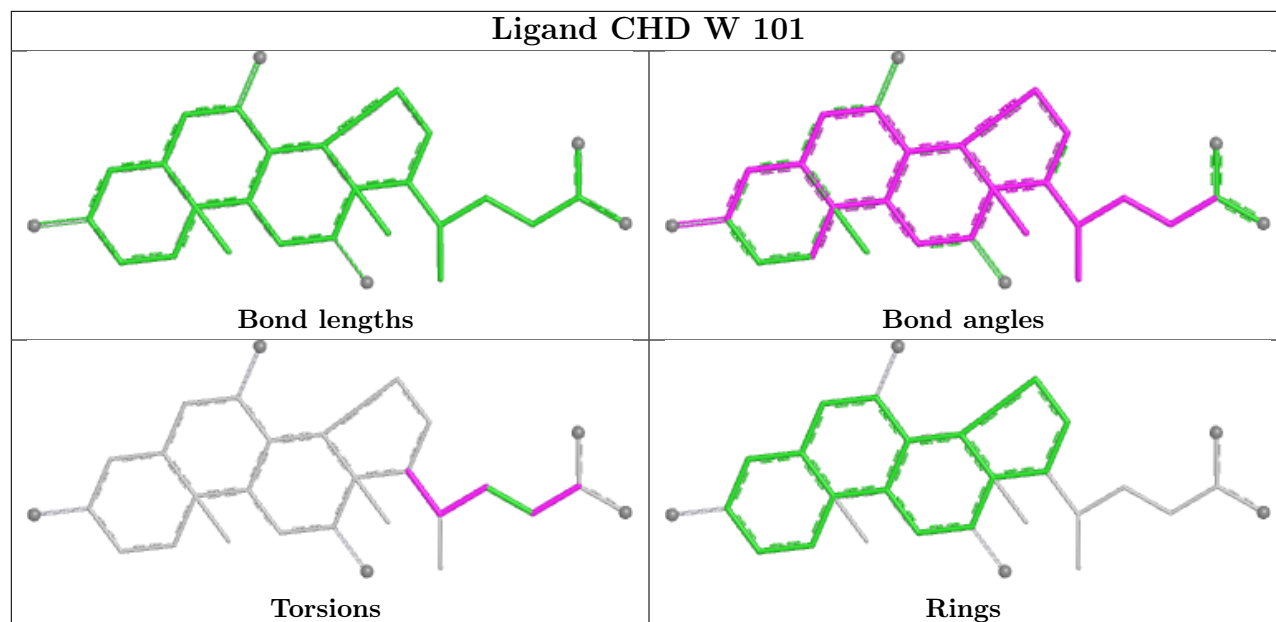
Ligand PEK T 101



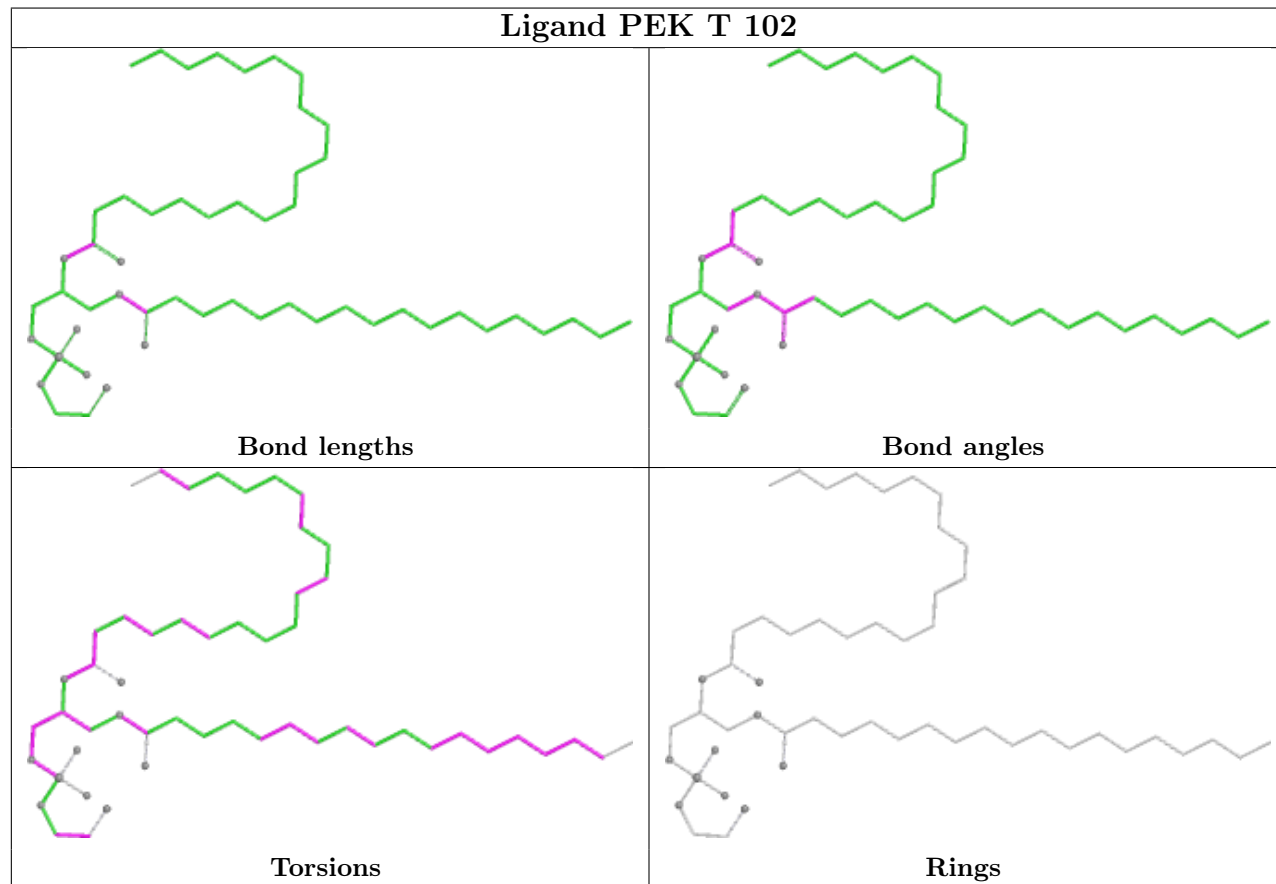


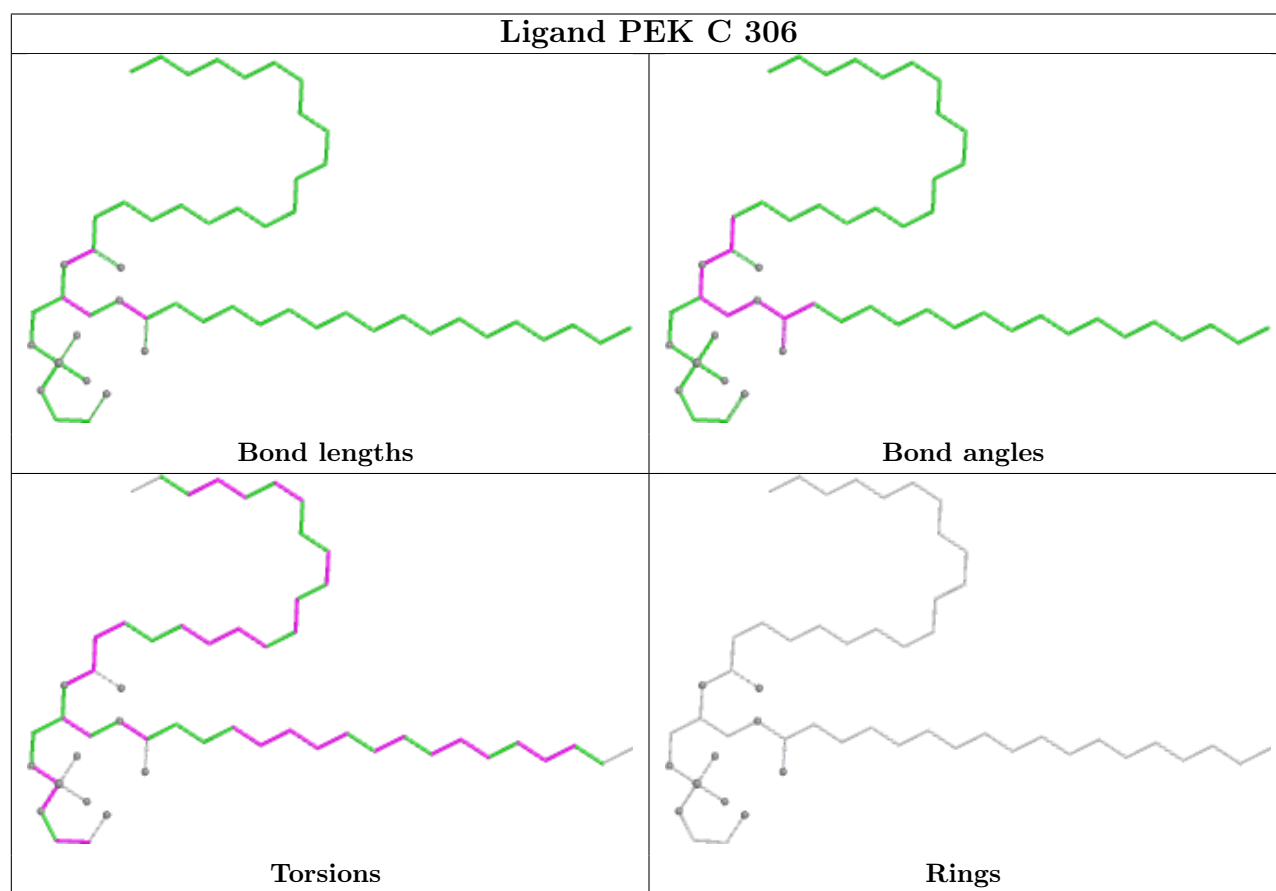
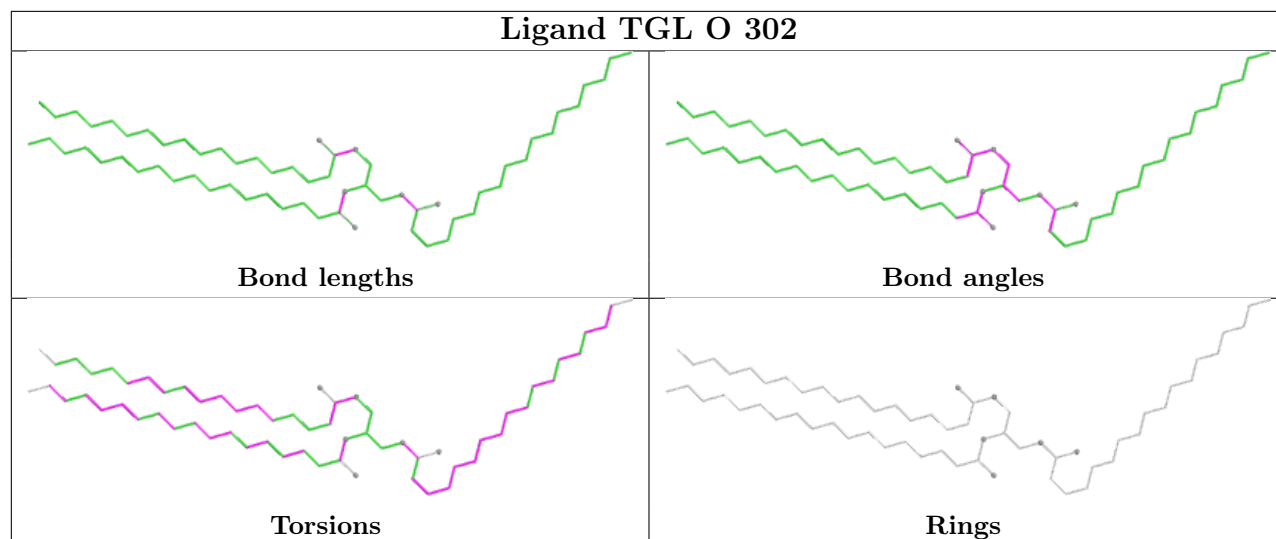


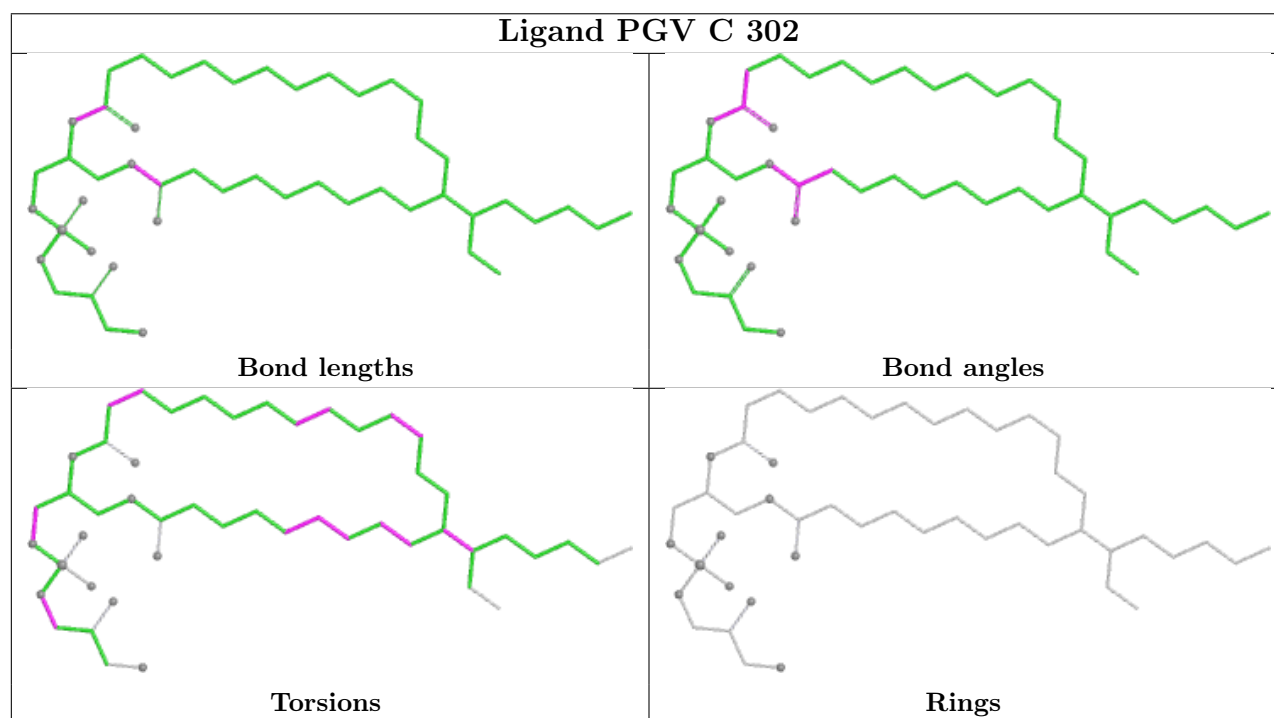
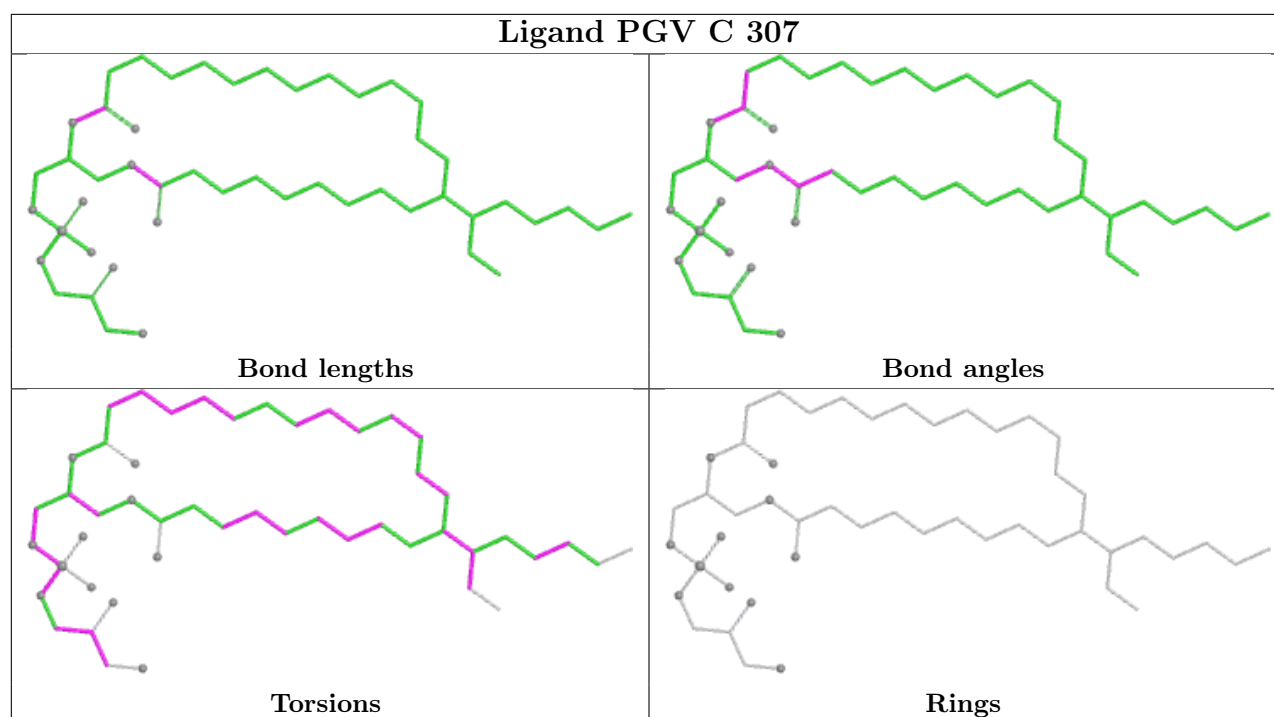
Ligand CHD W 101

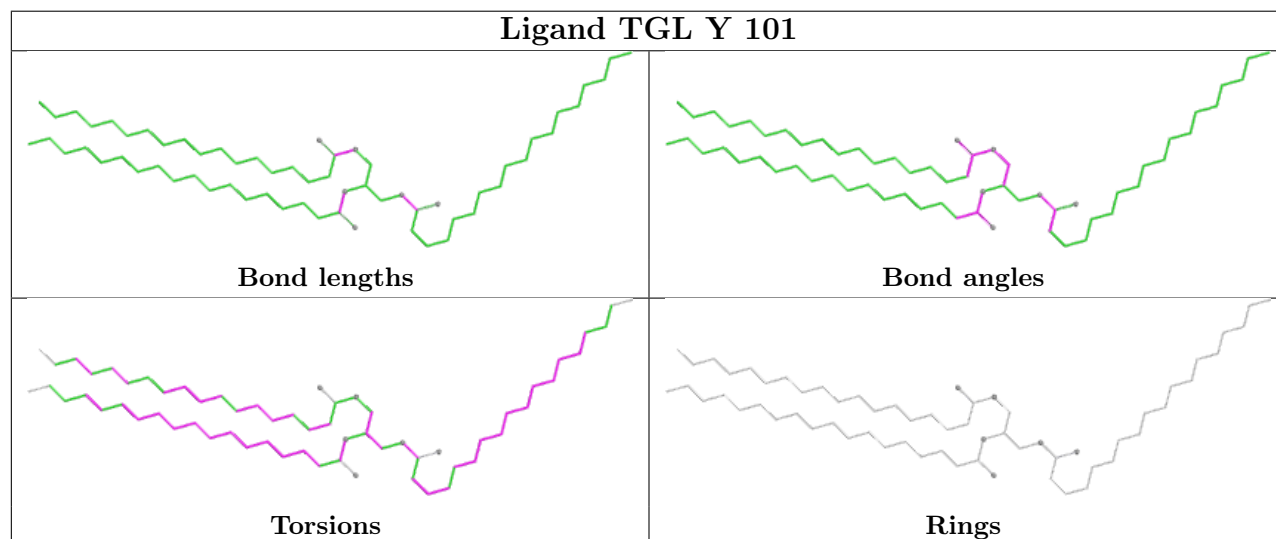
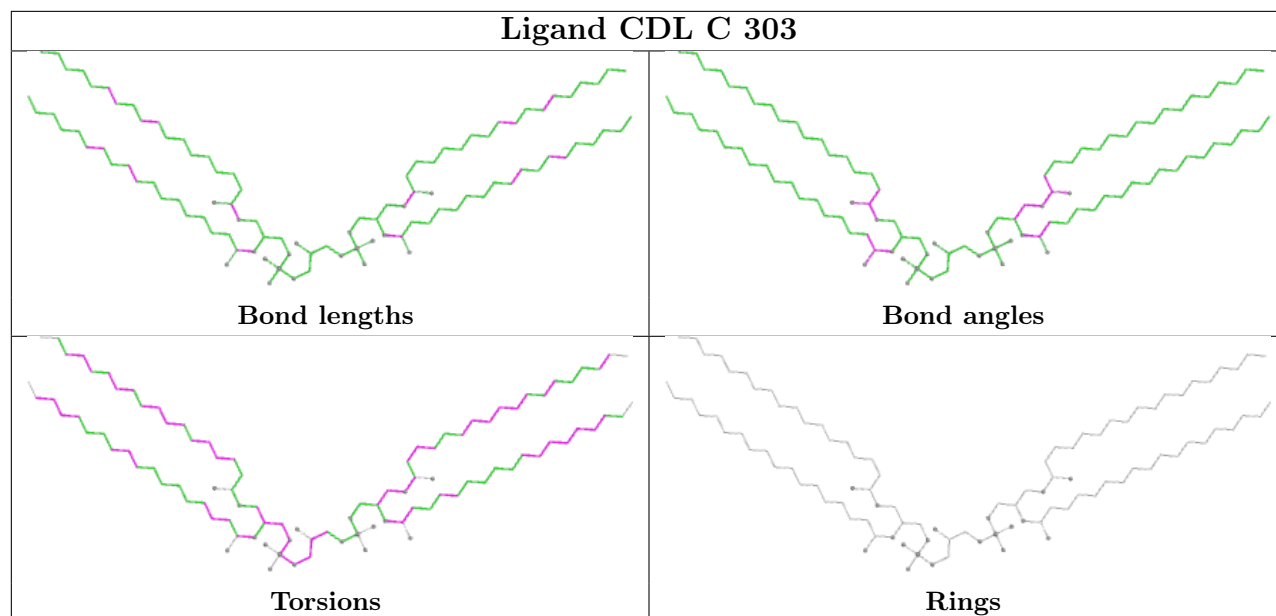


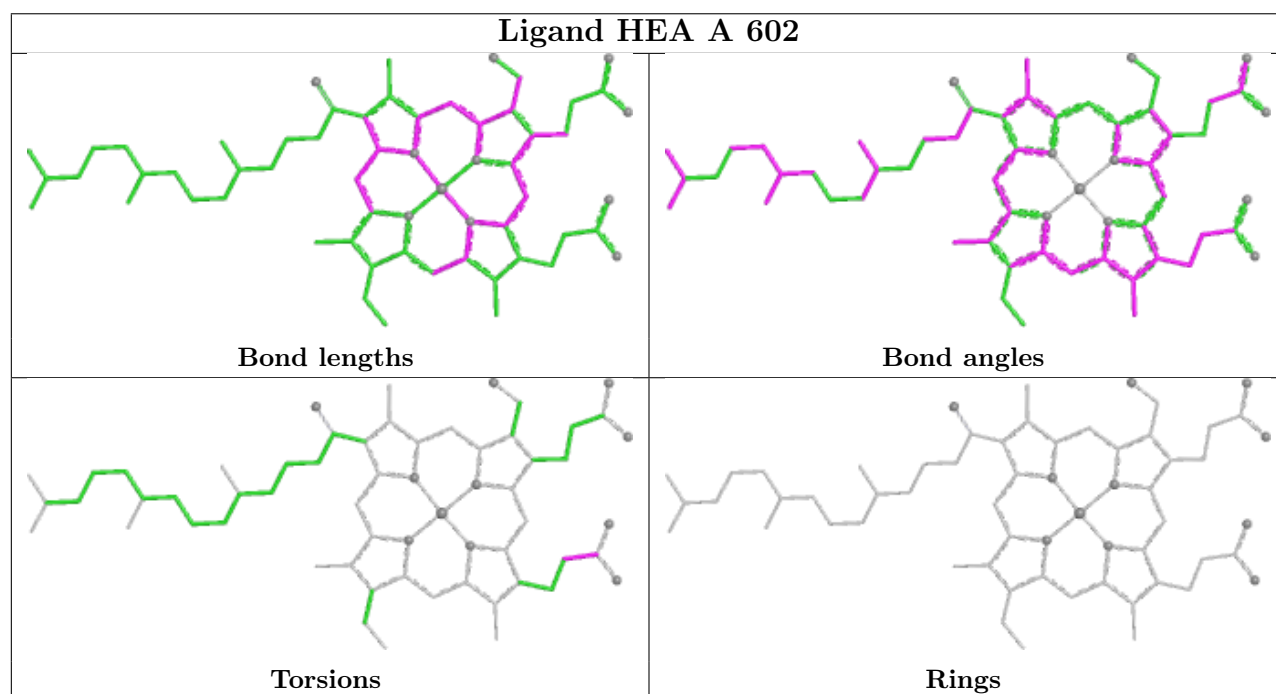
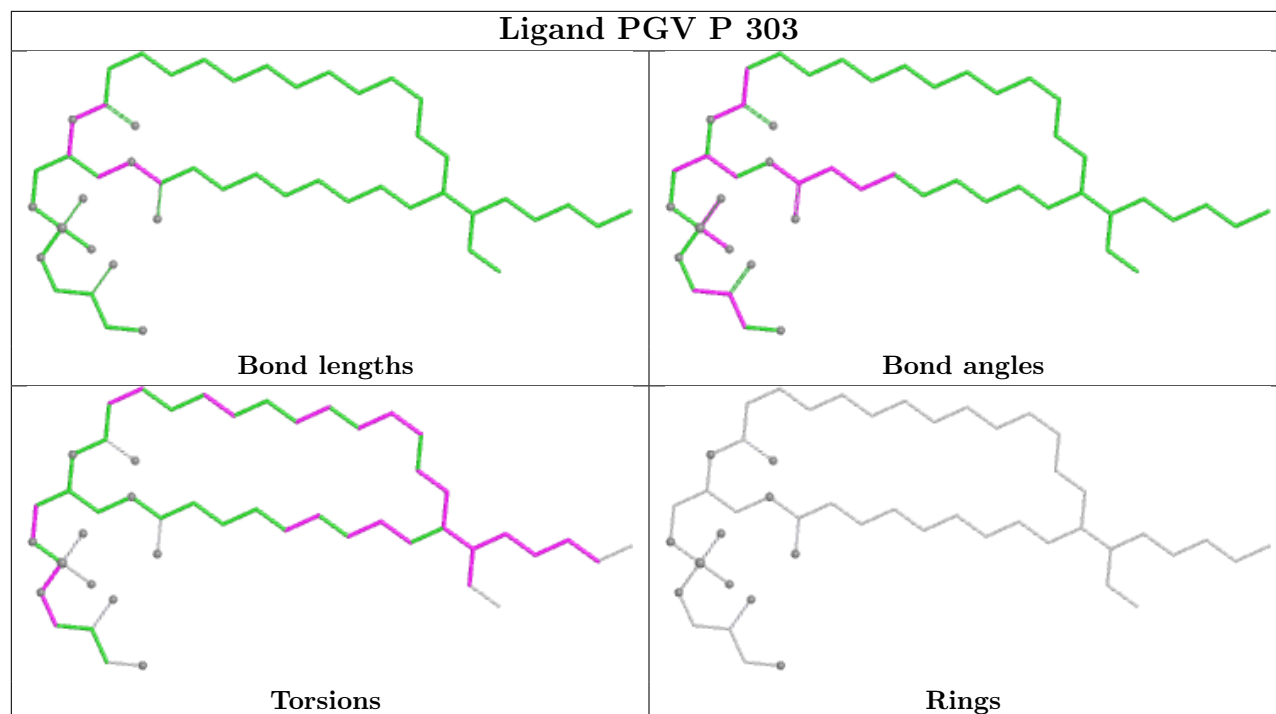
Ligand PEK T 102



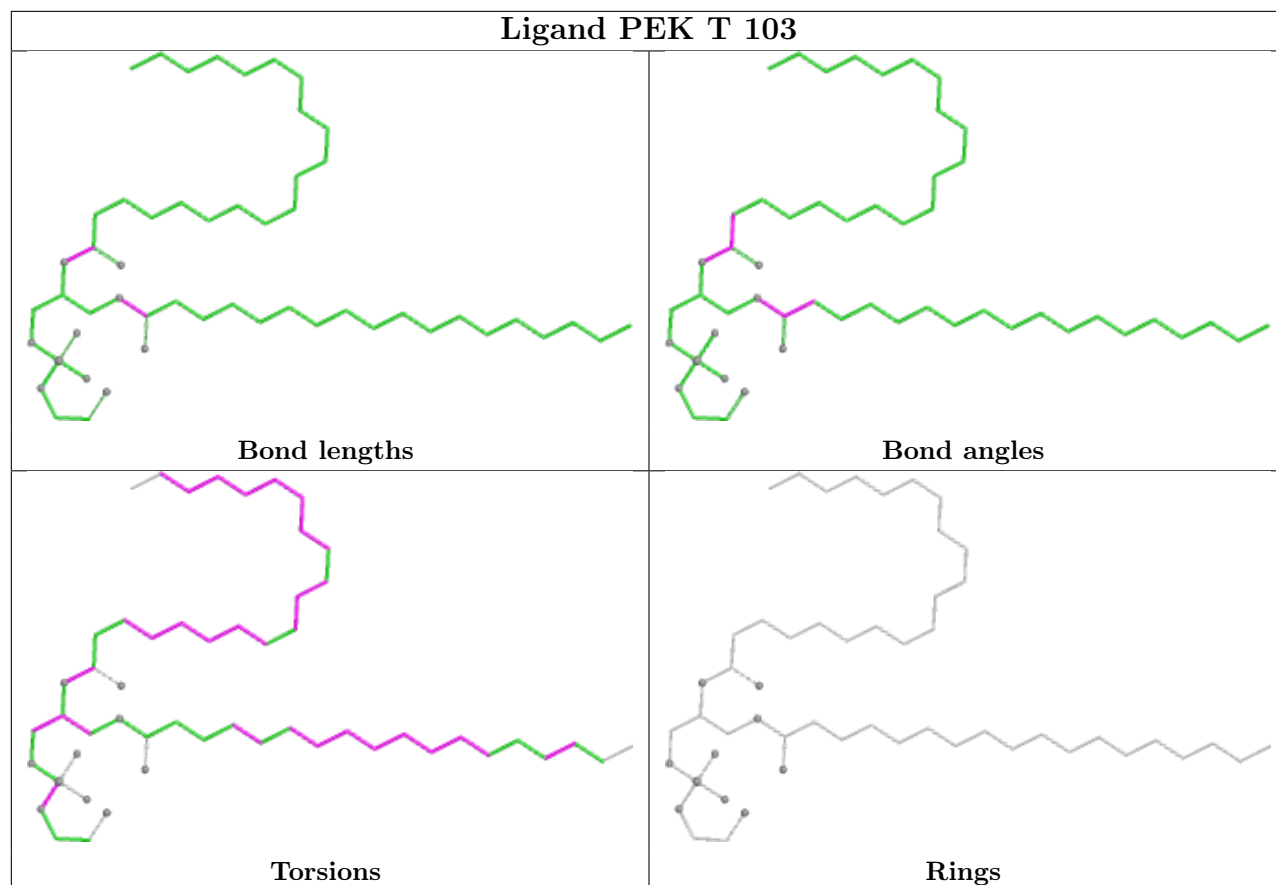




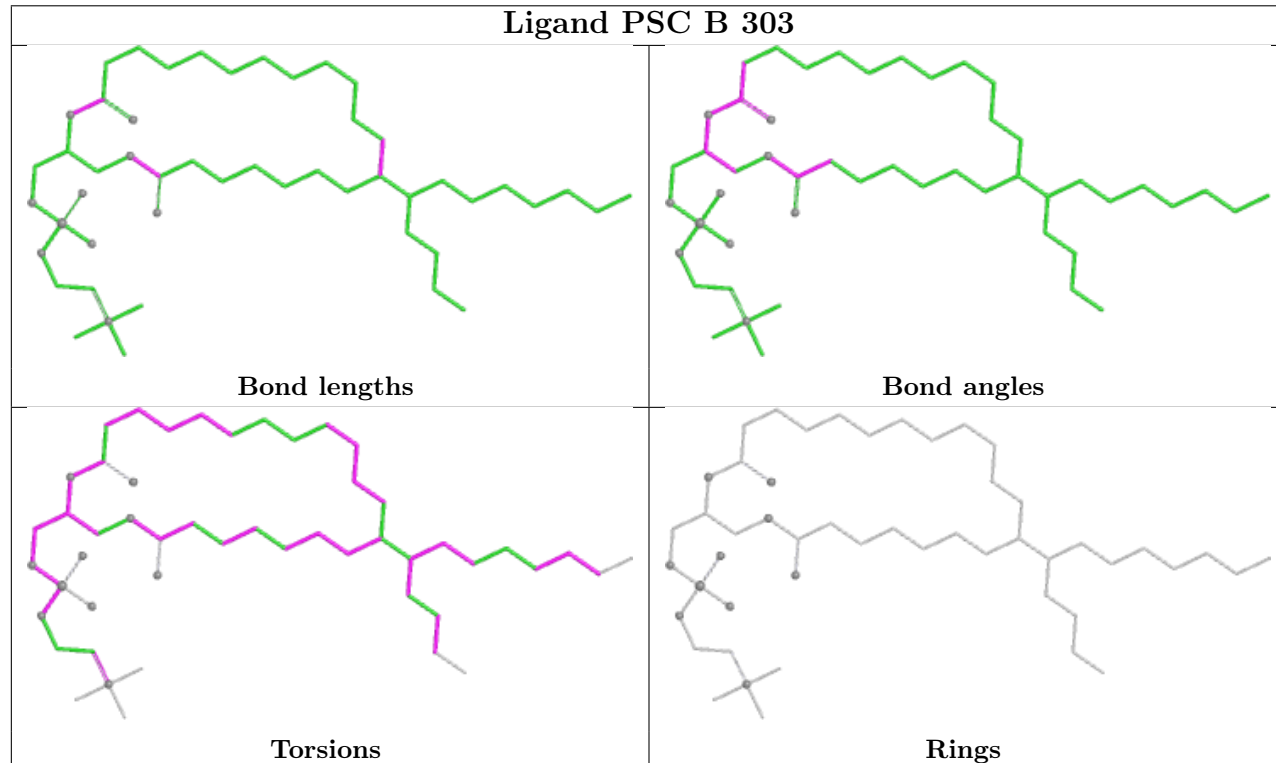


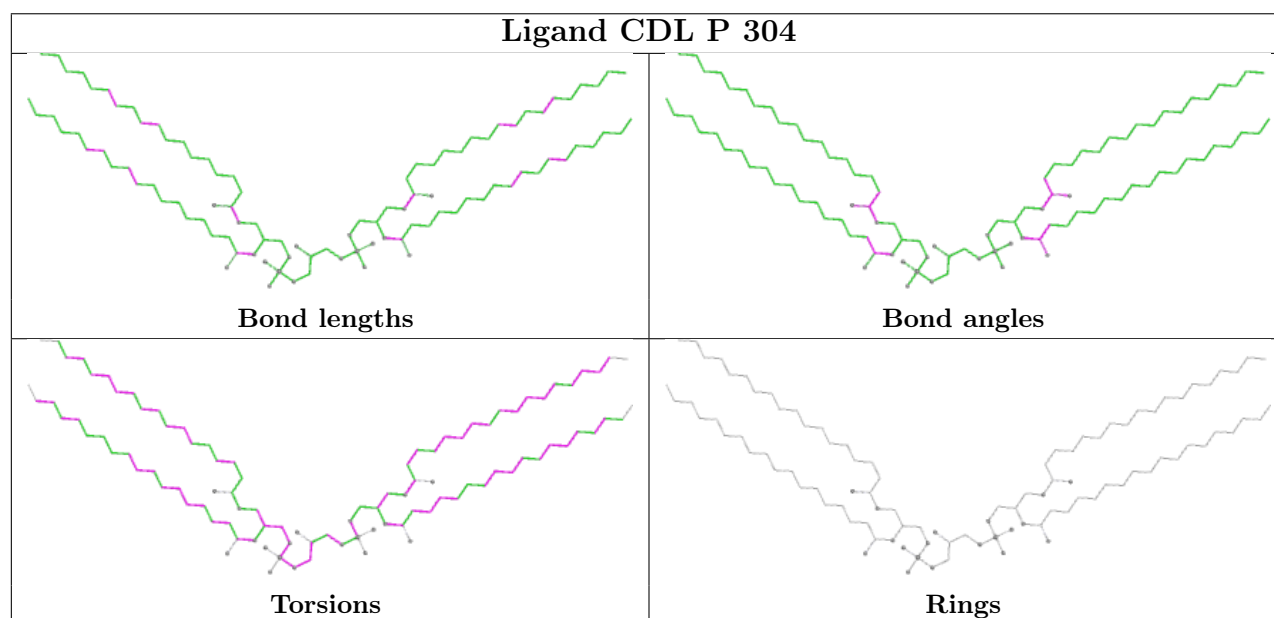
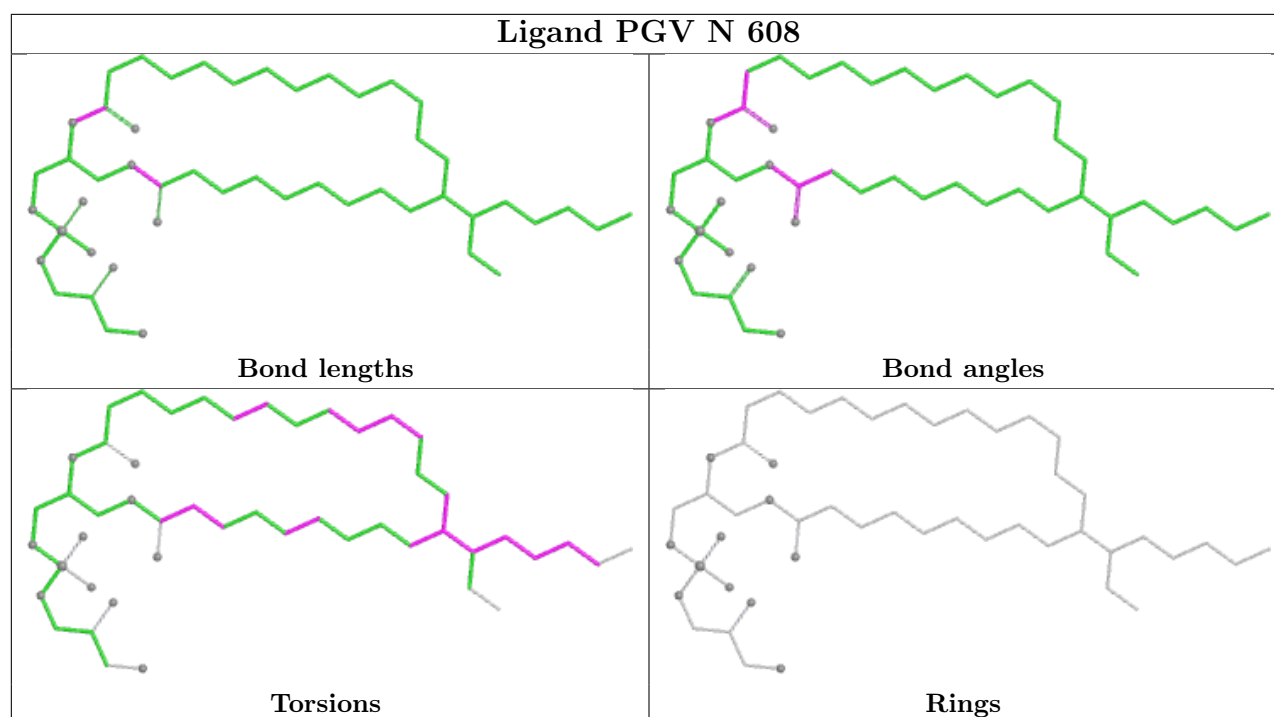


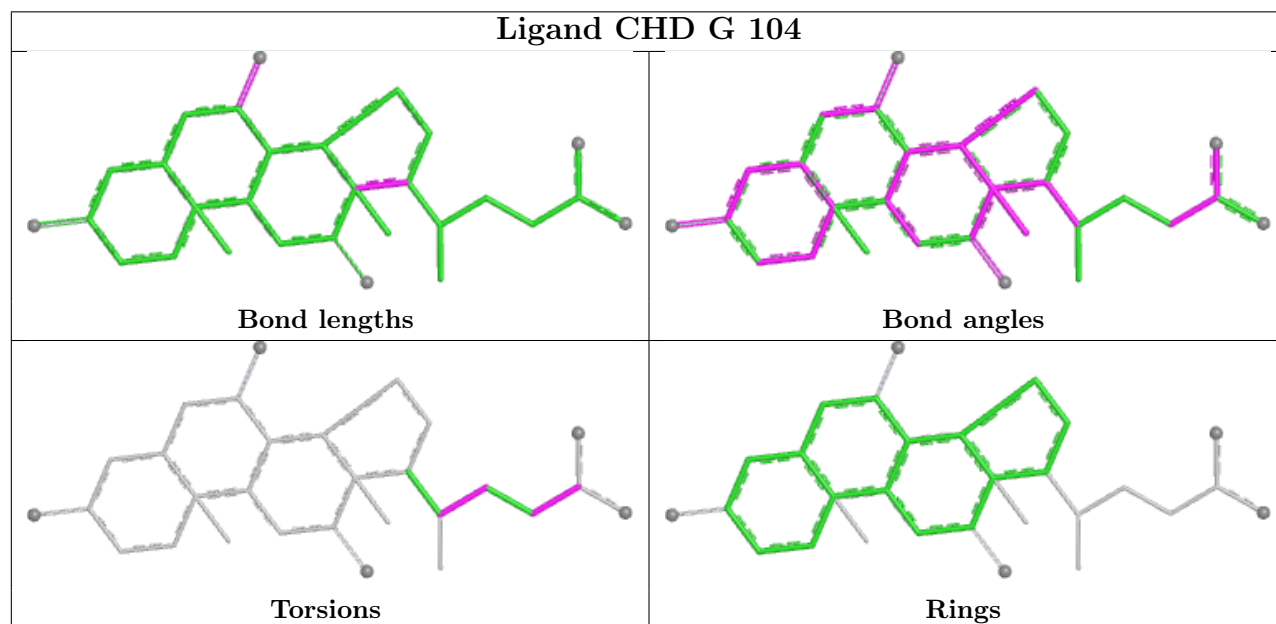
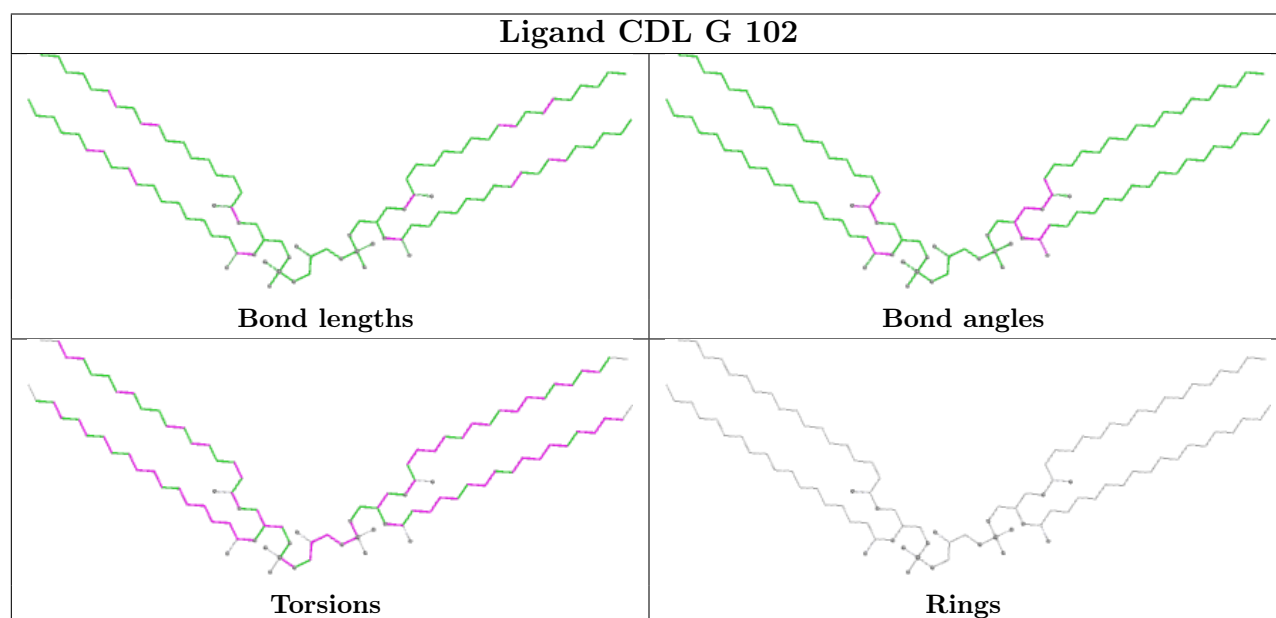
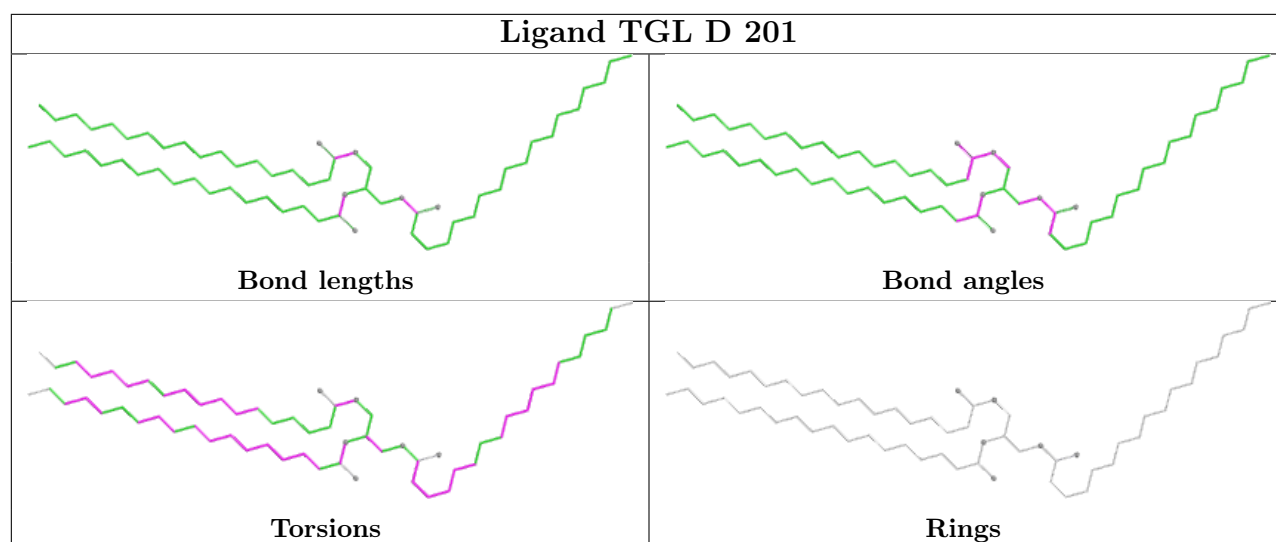
Ligand PEK T 103

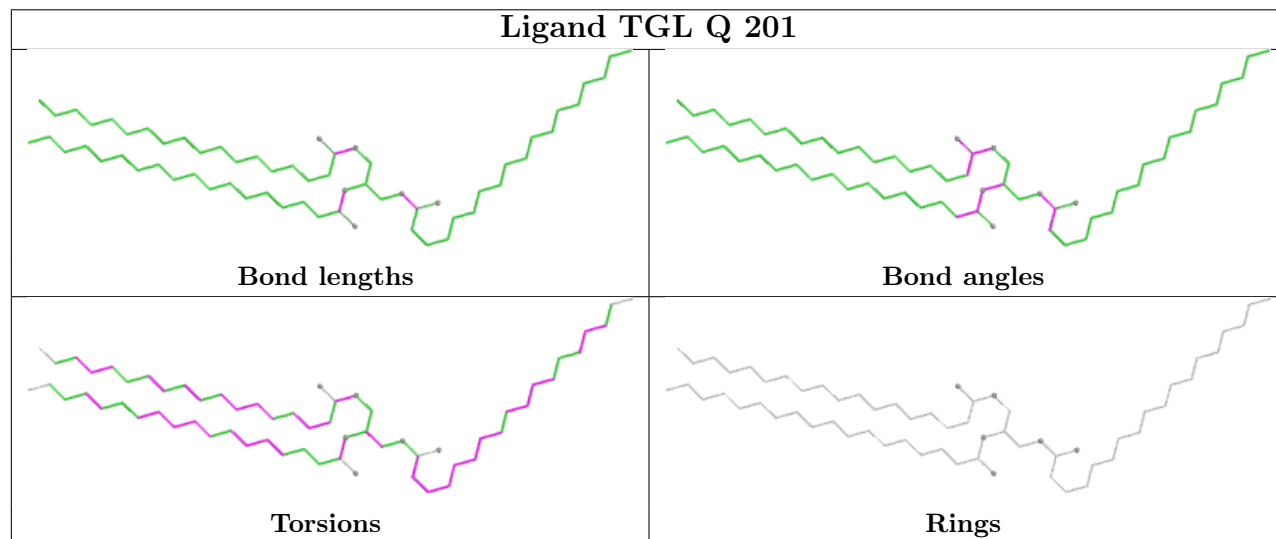
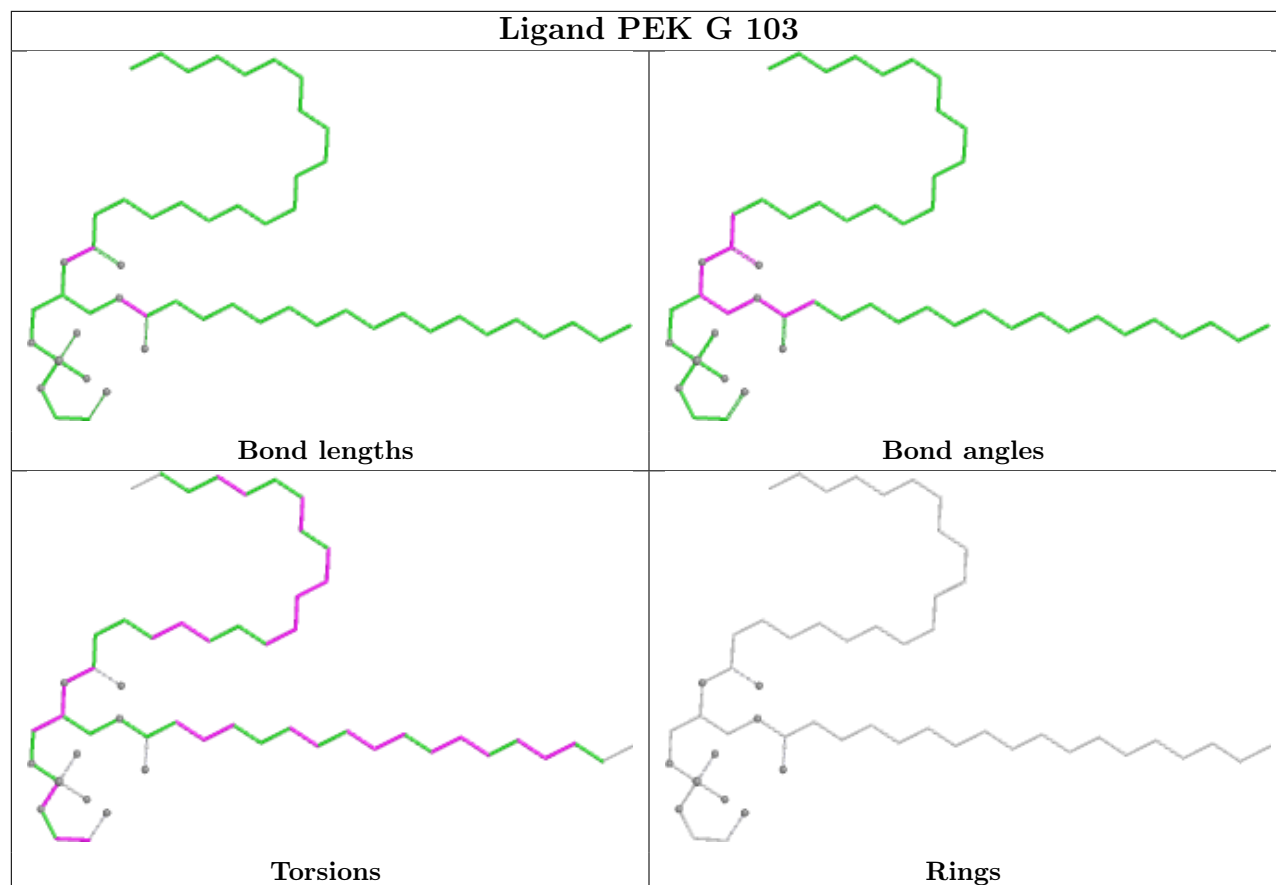


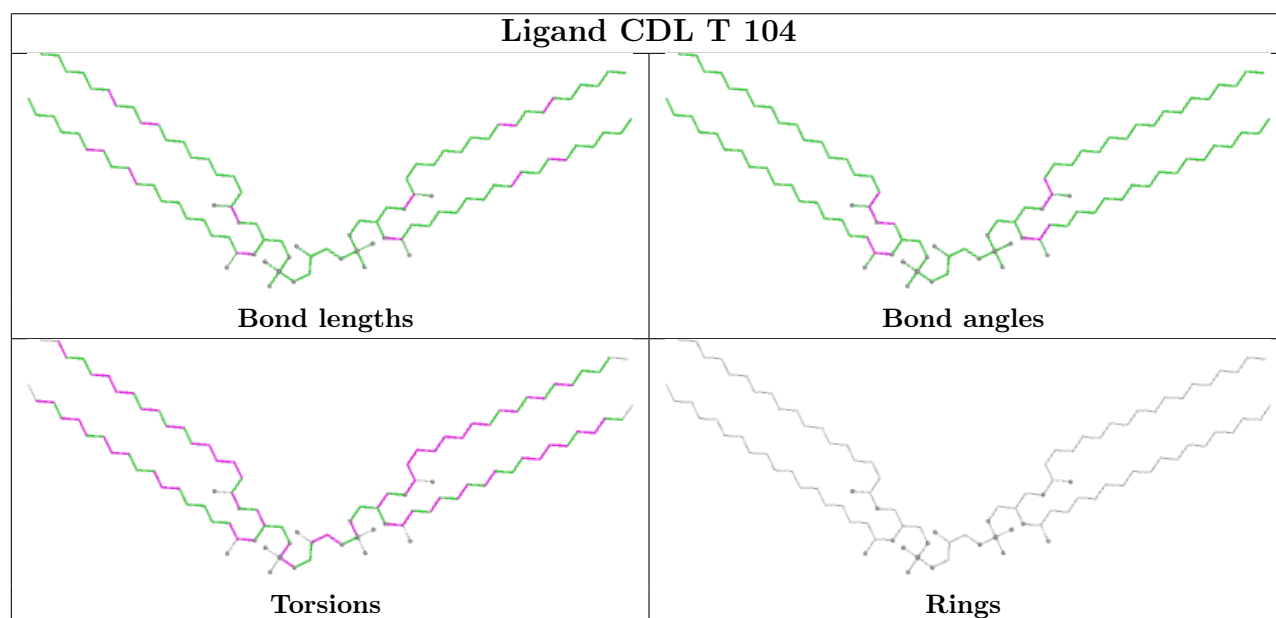
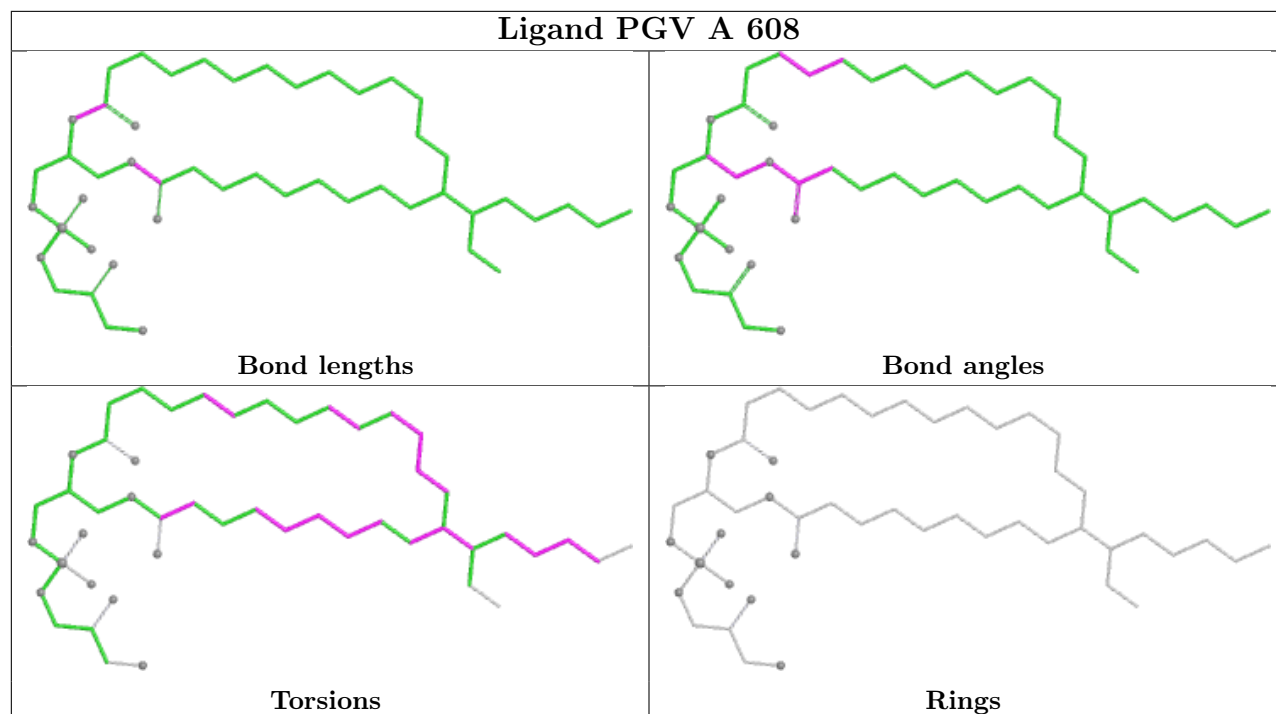
Ligand PSC B 303

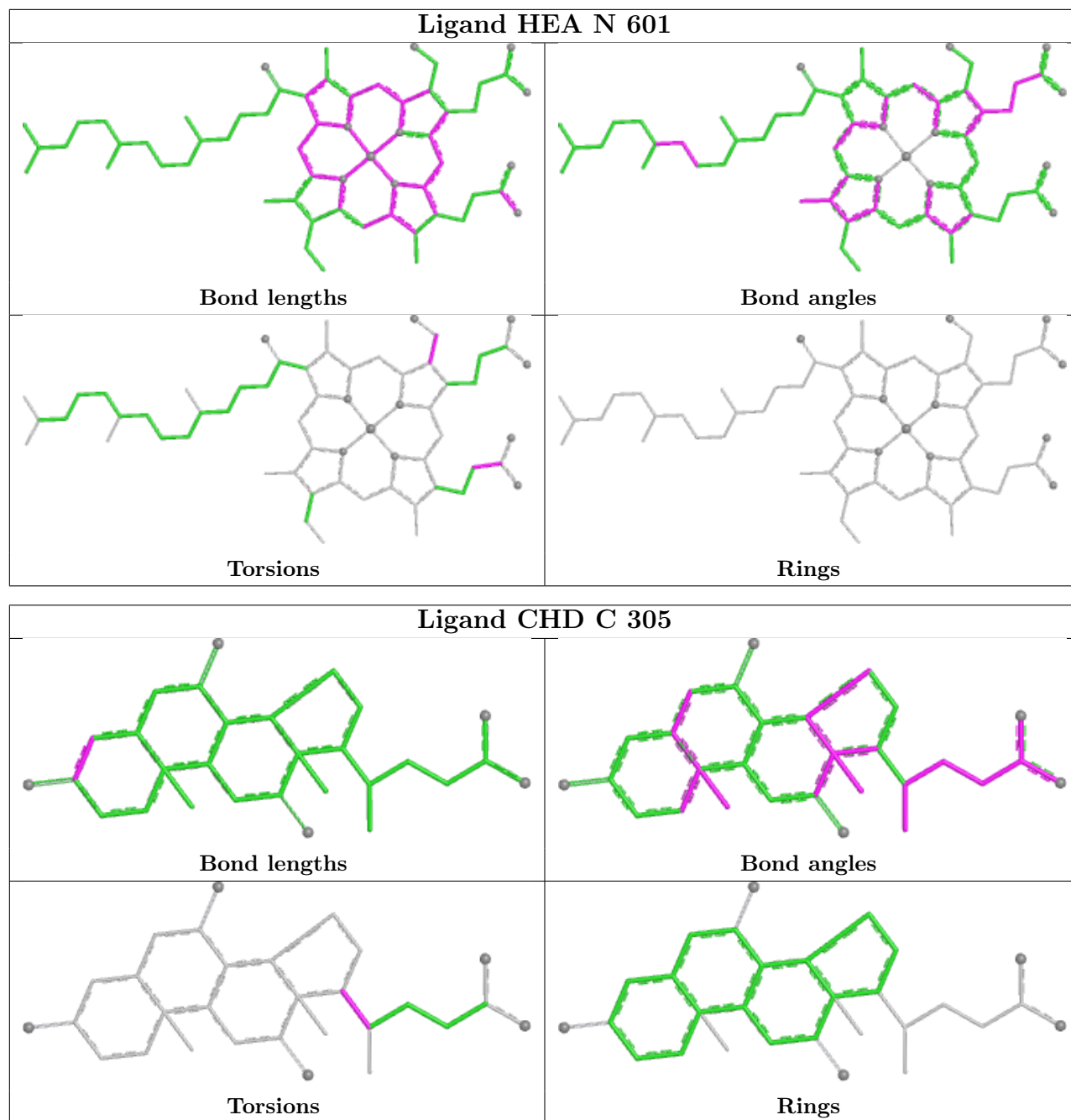


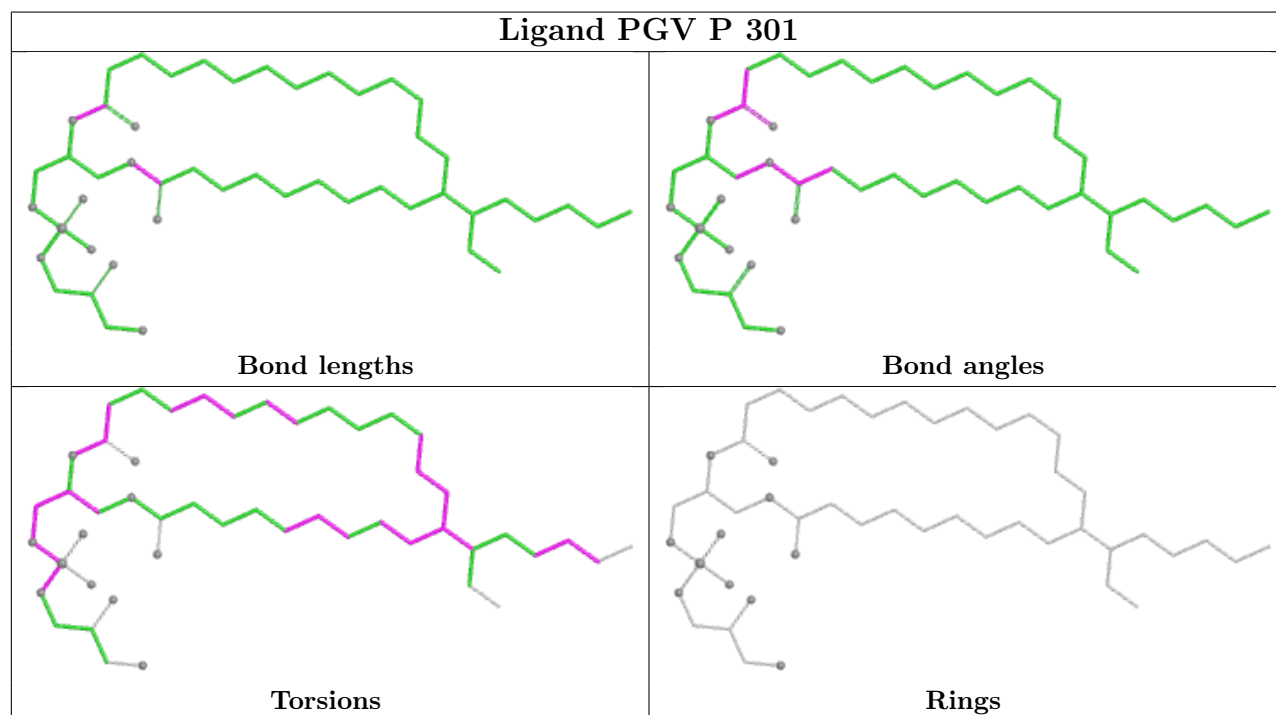
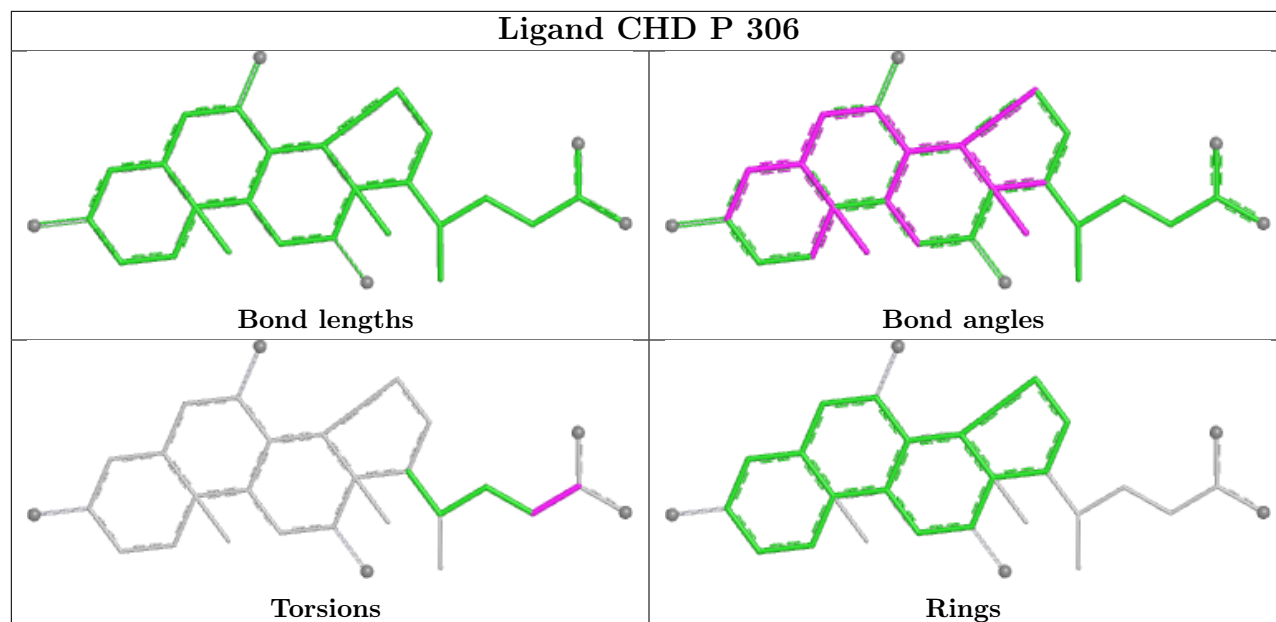


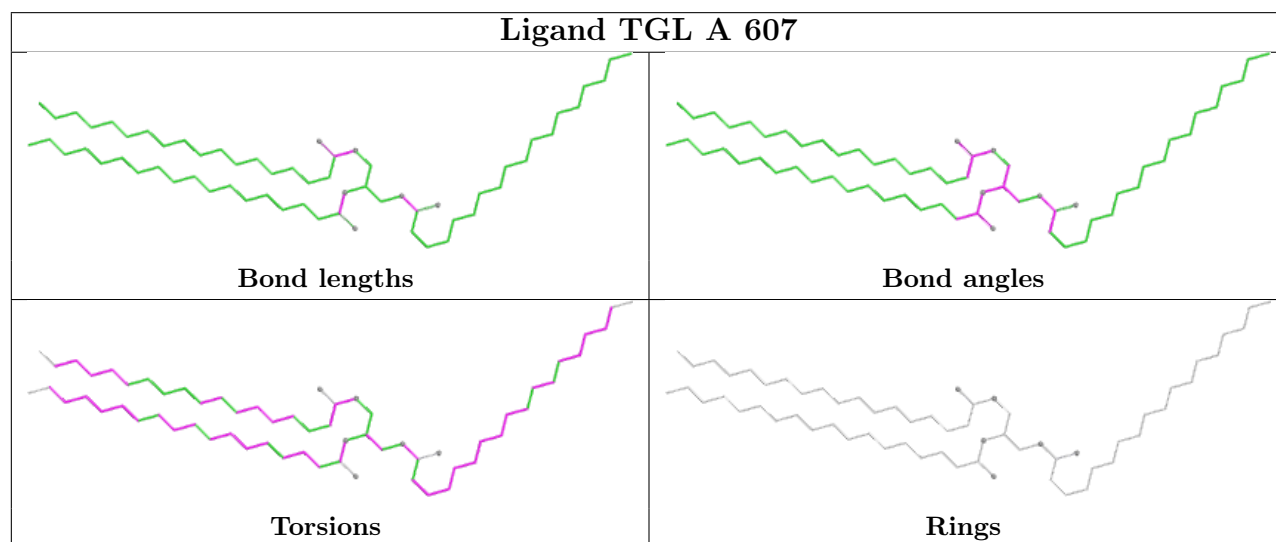
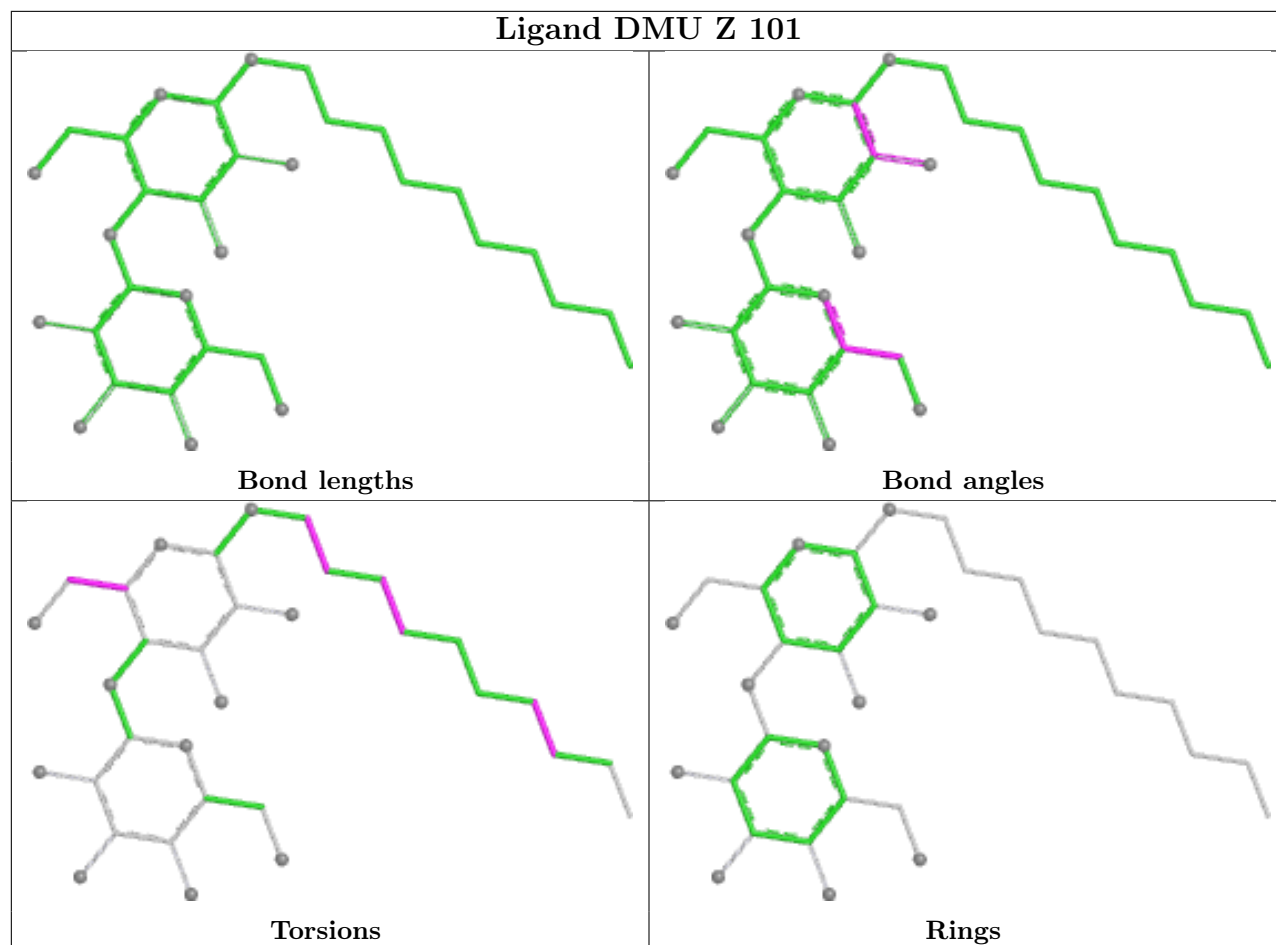


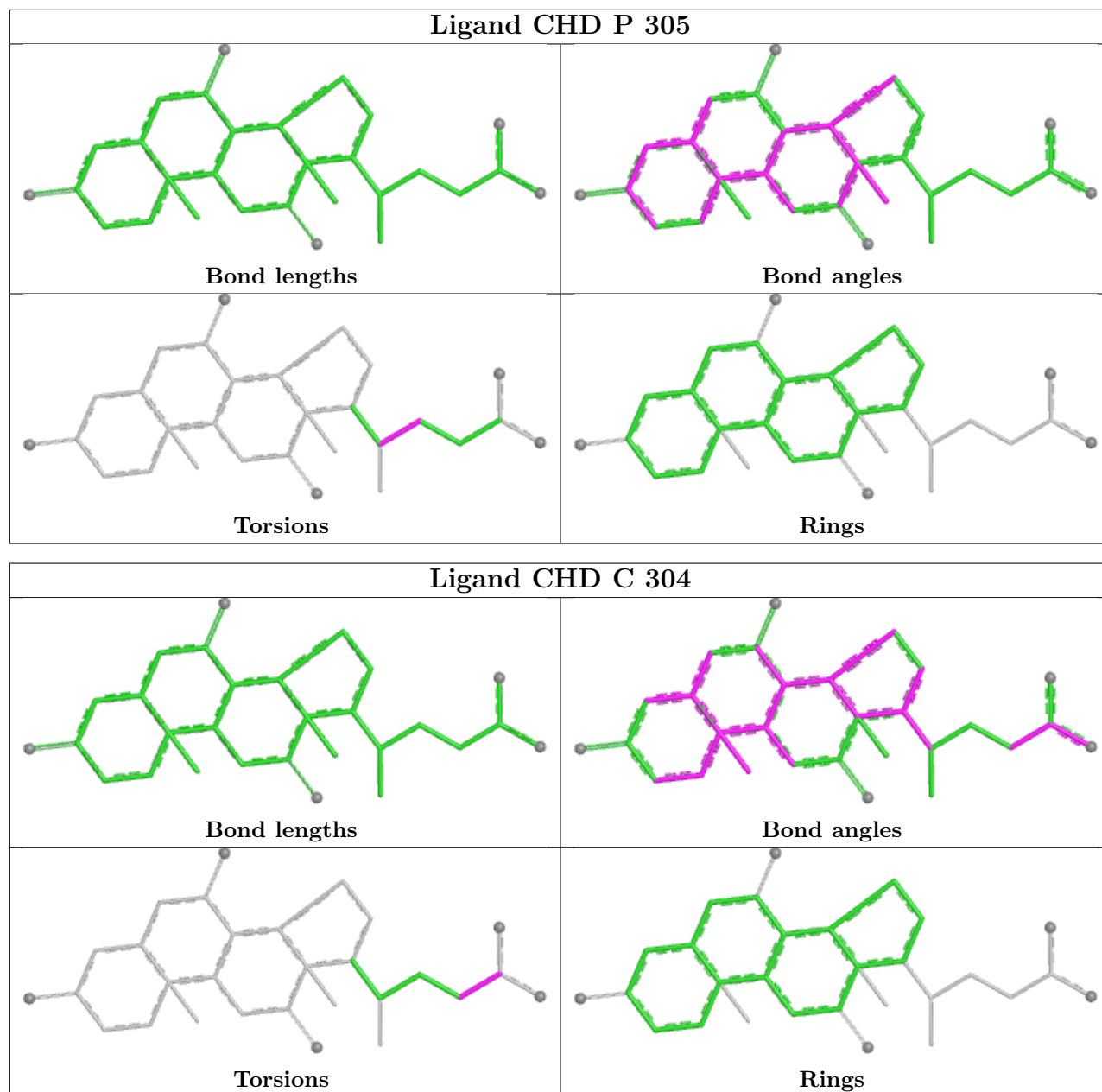


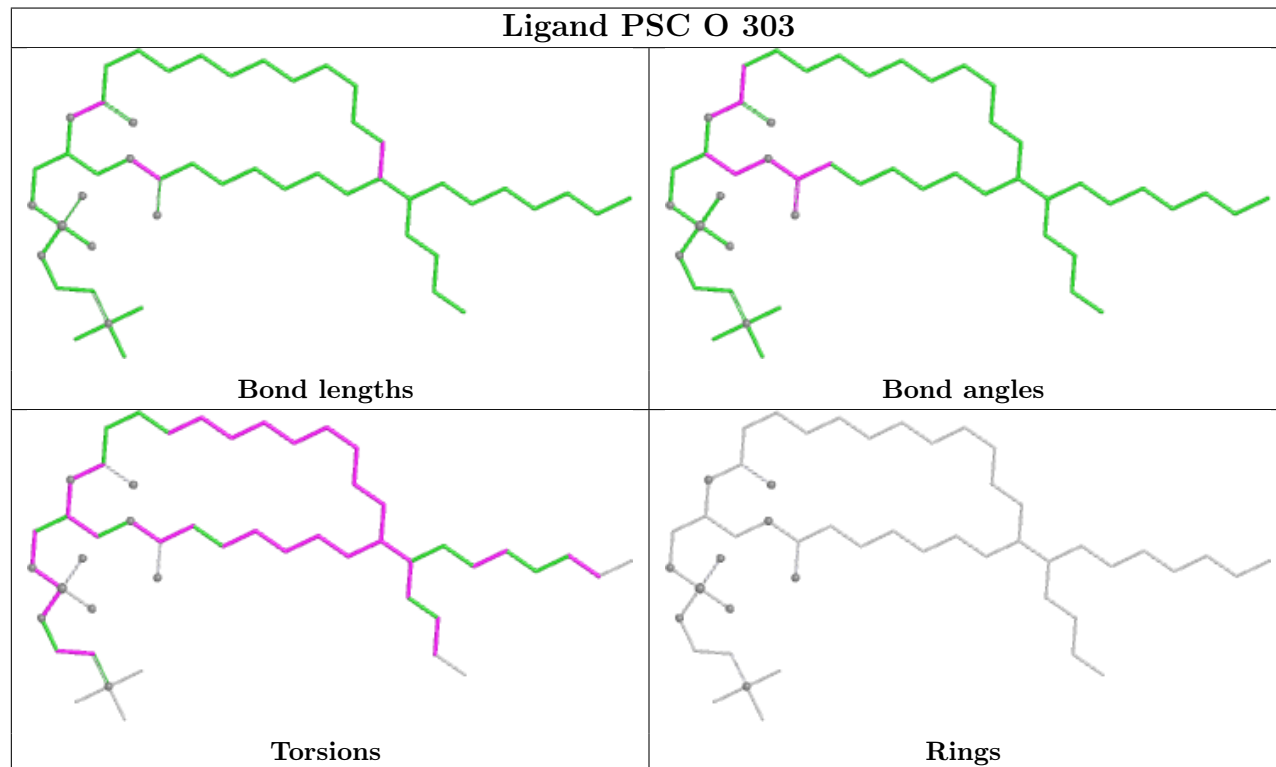
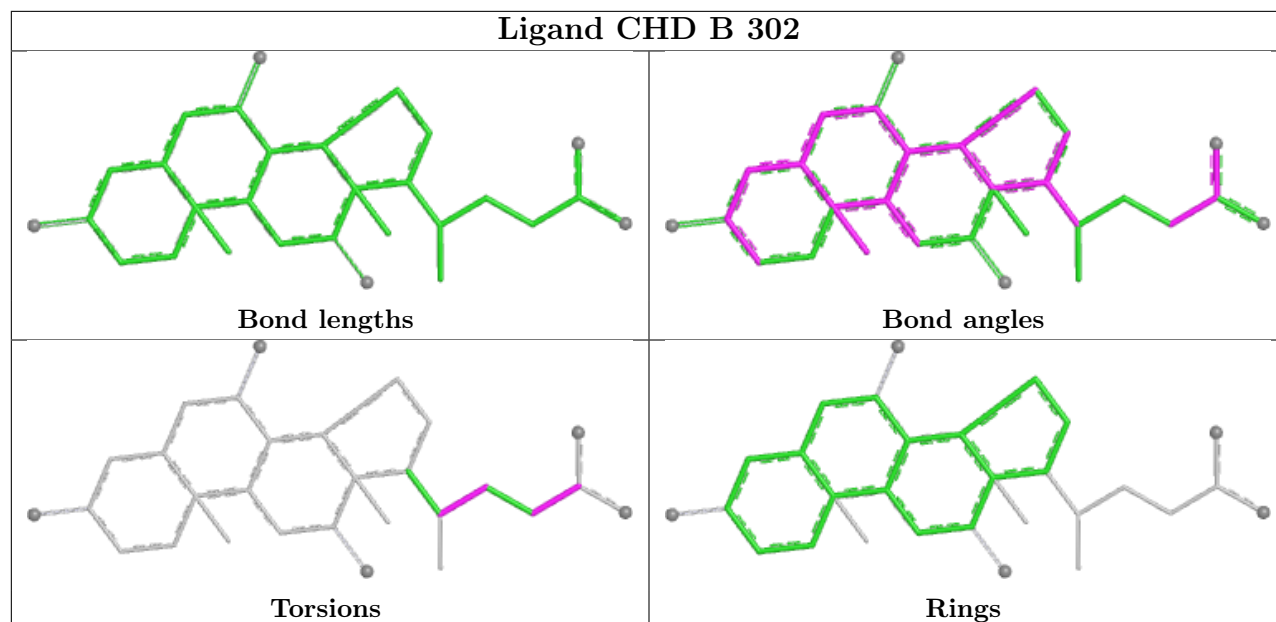


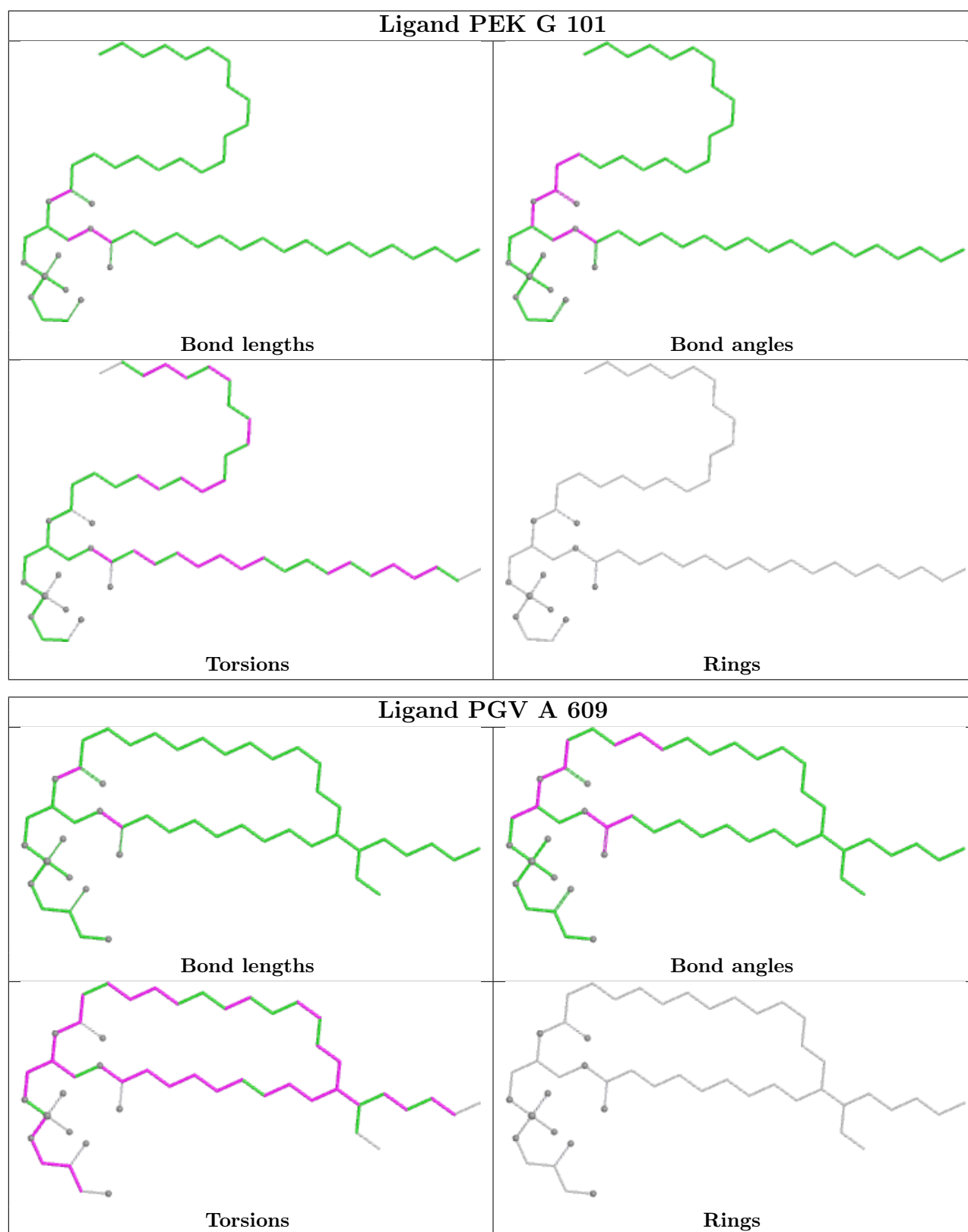












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	513/514 (99%)	-0.40	2 (0%) 88 87	23, 47, 59, 90	7 (1%)
1	N	513/514 (99%)	-0.21	10 (1%) 66 63	43, 55, 71, 103	0
2	B	226/227 (99%)	-0.33	1 (0%) 88 87	43, 54, 82, 118	0
2	O	226/227 (99%)	-0.09	5 (2%) 62 59	51, 64, 98, 140	0
3	C	259/261 (99%)	-0.35	1 (0%) 88 87	44, 51, 69, 123	0
3	P	259/261 (99%)	-0.09	6 (2%) 61 58	46, 56, 75, 129	0
4	D	144/147 (97%)	-0.17	1 (0%) 84 82	49, 59, 87, 126	0
4	Q	144/147 (97%)	0.32	11 (7%) 20 17	60, 81, 126, 157	0
5	E	105/109 (96%)	-0.26	1 (0%) 79 77	49, 59, 88, 147	0
5	R	105/109 (96%)	-0.18	3 (2%) 53 50	56, 73, 104, 142	0
6	F	98/98 (100%)	-0.21	5 (5%) 33 30	47, 61, 138, 155	0
6	S	98/98 (100%)	-0.00	3 (3%) 51 48	51, 70, 140, 158	0
7	G	83/85 (97%)	0.36	8 (9%) 13 11	49, 62, 150, 157	0
7	T	83/85 (97%)	0.27	9 (10%) 11 8	49, 68, 143, 154	0
8	H	79/85 (92%)	0.03	5 (6%) 26 23	49, 64, 129, 144	0
8	U	79/85 (92%)	-0.16	2 (2%) 58 55	56, 71, 143, 153	0
9	I	72/73 (98%)	-0.21	3 (4%) 40 37	44, 65, 96, 118	1 (1%)
9	V	72/73 (98%)	0.04	4 (5%) 30 27	46, 76, 105, 140	2 (2%)
10	J	58/59 (98%)	0.04	0 100 100	52, 64, 98, 150	0
10	W	58/59 (98%)	0.23	3 (5%) 33 29	59, 76, 110, 157	0
11	K	49/56 (87%)	0.01	2 (4%) 41 38	57, 67, 88, 105	0
11	X	49/56 (87%)	0.24	4 (8%) 17 15	69, 80, 109, 122	0
12	L	46/47 (97%)	-0.28	0 100 100	47, 55, 75, 109	0
12	Y	46/47 (97%)	0.27	4 (8%) 16 13	57, 72, 98, 140	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
13	M	43/46 (93%)	-0.03	1 (2%) 61 58	51, 57, 89, 130	0
13	Z	43/46 (93%)	0.07	0 100 100	66, 78, 113, 150	0
All	All	3550/3614 (98%)	-0.14	94 (2%) 57 54	23, 59, 103, 158	10 (0%)

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
7	G	5	LYS	9.0
7	G	4	ALA	7.9
7	G	3	ALA	7.1
6	S	96	LEU	6.9
9	V	37[A]	PHE	5.8
7	T	5	LYS	5.5
10	W	48	TYR	5.4
7	T	3	ALA	5.3
3	P	33	MET	5.2
7	G	2	SER	5.2
12	Y	37	PHE	5.1
7	G	6	GLY	5.0
6	S	97	ALA	4.9
3	P	37	PHE	4.8
6	S	1	ALA	4.8
9	I	37[A]	PHE	4.7
4	Q	140	TYR	4.7
1	N	2	PHE	4.6
2	O	90	ILE	4.2
4	Q	135	SER	4.1
6	F	1	ALA	4.0
9	V	38	ALA	4.0
6	F	65	ASP	3.9
7	T	2	SER	3.9
4	Q	5	VAL	3.8
1	A	296	GLY	3.6
4	Q	7	LYS	3.6
11	X	13	TYR	3.5
4	Q	6	VAL	3.5
7	T	4	ALA	3.5
11	X	40	TRP	3.5
10	W	49	CYS	3.3
4	D	4	SER	3.2
8	H	8	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
12	Y	19	TRP	3.2
4	Q	89	ILE	3.1
5	R	15	TRP	3.0
11	X	6	ALA	3.0
8	U	8	ILE	3.0
3	P	34	TRP	2.9
10	W	52	TRP	2.9
1	N	247	ILE	2.9
2	O	47	THR	2.9
2	O	43	SER	2.9
7	T	6	GLY	2.9
8	H	45	ALA	2.9
11	K	16	ALA	2.9
13	M	35	TYR	2.8
1	A	468	MET	2.8
3	P	39	SER	2.8
5	R	19	PHE	2.8
5	E	5	HIS	2.8
1	N	513	LEU	2.8
7	T	84	LYS	2.8
4	Q	125	ASP	2.7
8	H	24	ASN	2.7
8	H	7	LYS	2.7
7	T	1	ALA	2.7
7	T	7	ASP	2.6
1	N	484	THR	2.6
4	Q	93	ALA	2.6
4	Q	92	THR	2.6
7	T	9	GLY	2.6
1	N	222	PRO	2.6
12	Y	34	ALA	2.5
12	Y	3	TYR	2.5
1	N	397	PHE	2.4
1	N	57	VAL	2.4
1	N	470	PHE	2.4
11	X	17	VAL	2.4
2	O	211	LEU	2.4
5	R	104	LEU	2.4
1	N	472	ILE	2.3
9	I	25	PHE	2.3
7	G	10	GLY	2.3
9	I	21	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
6	F	96	LEU	2.2
11	K	19	ALA	2.2
7	G	36	TRP	2.2
1	N	381	LEU	2.2
3	P	122	HIS	2.2
6	F	64	GLU	2.2
4	Q	136	ALA	2.2
6	F	97	ALA	2.2
3	C	37	PHE	2.2
2	B	217	LYS	2.2
3	P	109	THR	2.2
9	V	34	PHE	2.2
4	Q	49	SER	2.2
8	H	84	LYS	2.1
7	G	9	GLY	2.1
2	O	136	LEU	2.0
9	V	31[A]	PHE	2.0
8	U	46	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
9	SAC	I	1	9/10	0.53	0.15	108,138,147,149	0
7	TPO	T	11	11/12	0.59	0.15	100,127,155,158	0
9	SAC	V	1	9/10	0.67	0.12	134,154,162,162	0
7	TPO	G	11	11/12	0.69	0.13	105,144,154,158	0
1	FME	N	1	10/11	0.83	0.20	77,90,144,145	0
1	FME	A	1	10/11	0.90	0.12	58,69,114,141	0
2	FME	B	1	10/11	0.98	0.09	51,56,62,76	0
2	FME	O	1	10/11	0.98	0.08	62,65,74,76	0

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
25	UNX	P	302	1/1	0.51	0.58	20,20,20,20	1
25	UNX	C	301	1/1	0.56	0.58	17,17,17,17	1
21	EDO	C	310	4/4	0.64	0.12	146,146,149,149	0
27	PEK	T	102	53/53	0.75	0.24	74,133,150,150	0
26	CDL	G	102	100/100	0.76	0.30	99,145,150,150	0
19	TGL	D	201	63/63	0.77	0.30	68,117,150,150	0
26	CDL	T	104	100/100	0.78	0.29	83,136,150,150	0
26	CDL	P	304	100/100	0.79	0.22	70,139,150,150	0
23	CHD	W	101	29/29	0.80	0.13	108,149,150,150	0
27	PEK	G	103	53/53	0.81	0.19	84,136,150,150	0
19	TGL	A	607	63/63	0.82	0.21	50,96,139,151	0
19	TGL	Y	101	63/63	0.82	0.27	78,127,150,150	0
20	PGV	P	301	51/51	0.82	0.17	73,119,150,150	0
21	EDO	N	610	4/4	0.83	0.17	125,132,135,138	0
27	PEK	C	306	53/53	0.83	0.23	68,138,150,150	0
19	TGL	O	302	63/63	0.83	0.22	67,115,150,150	0
20	PGV	A	609	51/51	0.83	0.26	62,112,152,160	0
20	PGV	N	607	51/51	0.84	0.24	66,140,150,150	0
21	EDO	F	103	4/4	0.84	0.27	71,125,128,147	0
23	CHD	J	101	29/29	0.85	0.15	90,142,150,150	0
19	TGL	Q	201	63/63	0.85	0.25	76,124,150,150	0
24	PSC	B	303	52/52	0.86	0.19	73,142,150,150	0
23	CHD	P	305	29/29	0.86	0.15	82,131,147,148	0
20	PGV	C	307	51/51	0.86	0.19	61,126,150,150	0
19	TGL	L	101	63/63	0.87	0.25	67,106,149,150	0
26	CDL	C	303	100/100	0.87	0.15	56,121,153,157	0
21	EDO	A	611	4/4	0.87	0.21	81,82,106,109	0
21	EDO	K	102	4/4	0.87	0.15	89,96,101,112	0
27	PEK	T	103	53/53	0.88	0.21	67,132,150,150	0
21	EDO	H	101	4/4	0.89	0.25	111,112,118,127	0
21	EDO	C	308	4/4	0.90	0.13	89,93,96,106	0
23	CHD	C	304	29/29	0.90	0.09	89,104,112,113	0
21	EDO	A	613	4/4	0.91	0.13	96,116,120,122	0
21	EDO	C	311	4/4	0.91	0.17	60,77,78,84	0
21	EDO	P	307	4/4	0.91	0.16	92,96,107,115	0
29	DMU	Z	101	33/33	0.91	0.13	95,106,138,147	0
21	EDO	P	308	4/4	0.92	0.13	89,95,98,99	0

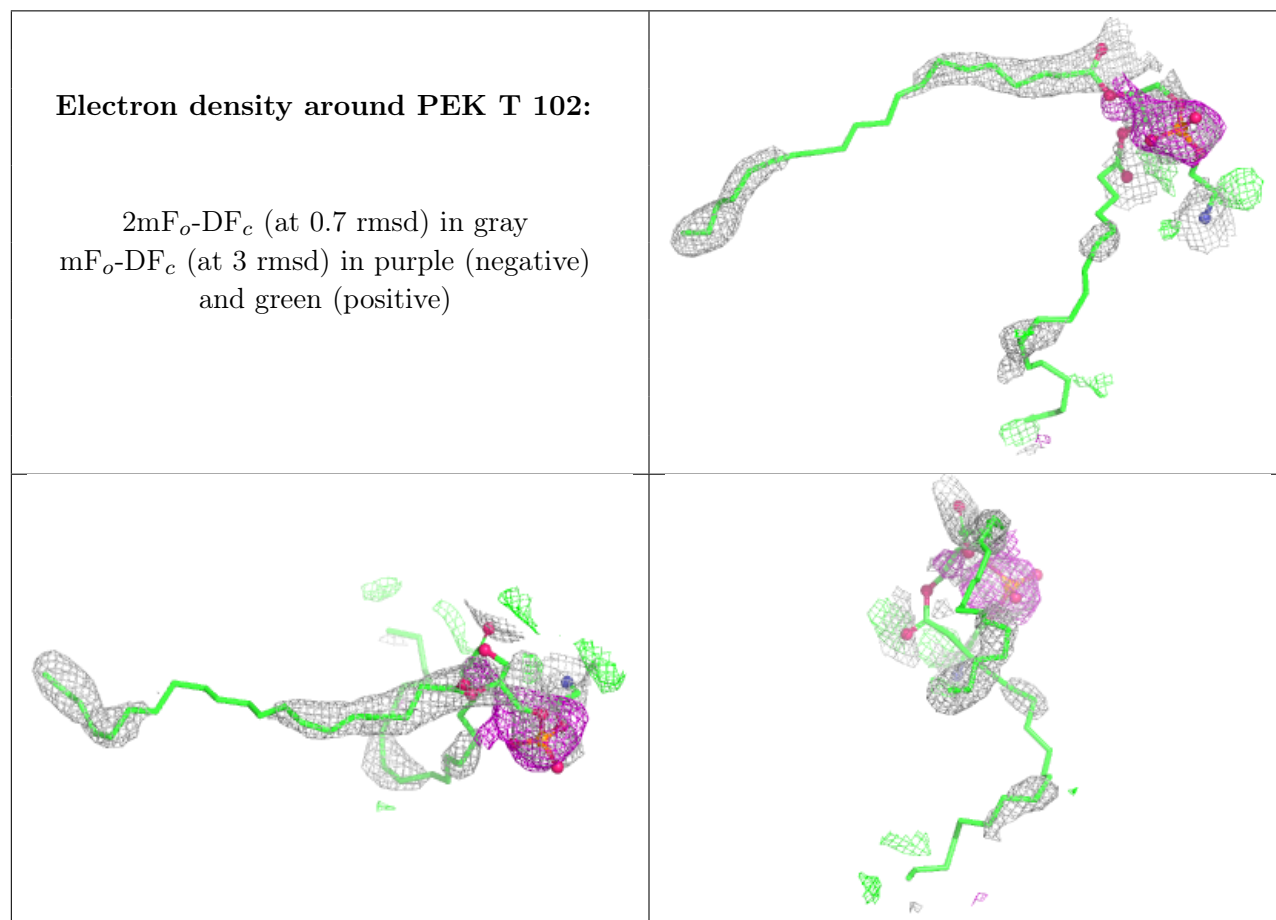
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
24	PSC	O	303	52/52	0.93	0.16	72,134,150,150	0
21	EDO	L	102	4/4	0.93	0.11	83,86,90,95	0
21	EDO	B	305	4/4	0.93	0.17	84,86,91,94	0
21	EDO	C	309	4/4	0.93	0.11	80,86,89,95	0
21	EDO	F	102	4/4	0.93	0.12	82,82,101,101	0
21	EDO	T	105	4/4	0.93	0.16	121,130,135,136	0
21	EDO	K	101	4/4	0.94	0.09	87,91,96,105	0
29	DMU	M	101	33/33	0.94	0.09	61,78,111,115	0
23	CHD	P	306	29/29	0.94	0.09	52,60,64,66	0
18	CMO	N	606	2/2	0.95	0.20	47,47,47,50	2
21	EDO	O	304	4/4	0.95	0.28	68,70,73,77	0
21	EDO	A	610	4/4	0.96	0.14	54,63,68,71	0
20	PGV	P	303	51/51	0.96	0.12	53,65,143,148	0
27	PEK	T	101	53/53	0.96	0.12	62,77,149,150	0
21	EDO	G	105	4/4	0.96	0.09	74,78,88,97	0
21	EDO	N	609	4/4	0.96	0.11	88,91,100,108	0
21	EDO	A	612	4/4	0.96	0.12	62,76,80,93	0
23	CHD	C	305	29/29	0.96	0.07	48,56,63,68	0
20	PGV	C	302	51/51	0.97	0.10	46,57,107,115	0
18	CMO	A	606	2/2	0.97	0.39	33,33,33,35	2
27	PEK	G	101	53/53	0.97	0.10	50,67,123,137	0
14	HEA	A	602	60/60	0.97	0.08	37,44,53,64	0
20	PGV	N	608	51/51	0.97	0.11	50,71,101,113	0
23	CHD	G	104	29/29	0.97	0.07	42,51,59,66	0
21	EDO	B	304	4/4	0.97	0.17	64,66,67,75	0
20	PGV	A	608	51/51	0.97	0.09	43,65,86,105	0
17	NA	N	605	1/1	0.97	0.04	65,65,65,65	0
23	CHD	B	302	29/29	0.98	0.05	49,57,63,73	0
14	HEA	N	601	60/60	0.98	0.08	48,57,76,84	0
14	HEA	N	602	60/60	0.98	0.07	40,49,58,65	0
14	HEA	A	601	60/60	0.98	0.07	38,44,59,72	0
16	MG	N	604	1/1	0.99	0.02	52,52,52,52	0
28	ZN	S	101	1/1	0.99	0.03	64,64,64,64	0
17	NA	A	605	1/1	0.99	0.04	53,53,53,53	0
22	CUA	O	301	2/2	0.99	0.02	56,56,56,59	0
16	MG	A	604	1/1	1.00	0.03	41,41,41,41	0
28	ZN	F	101	1/1	1.00	0.03	57,57,57,57	0
22	CUA	B	301	2/2	1.00	0.03	48,48,48,50	0
15	CU	A	603	1/1	1.00	0.01	45,45,45,45	0
15	CU	N	603	1/1	1.00	0.01	51,51,51,51	0

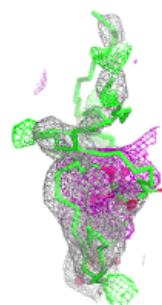
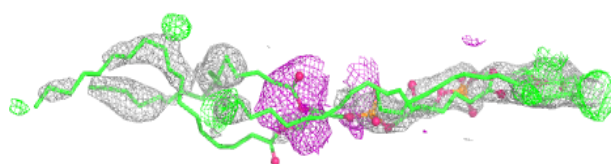
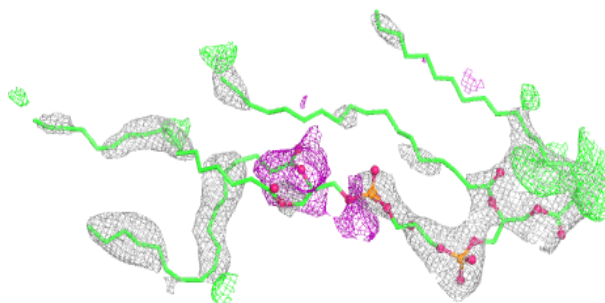
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

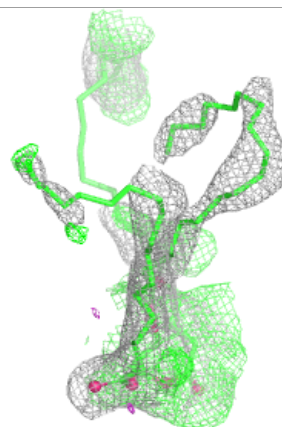
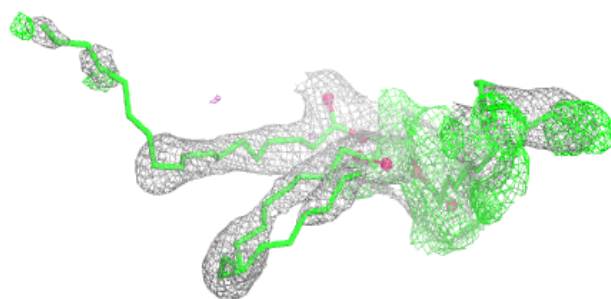
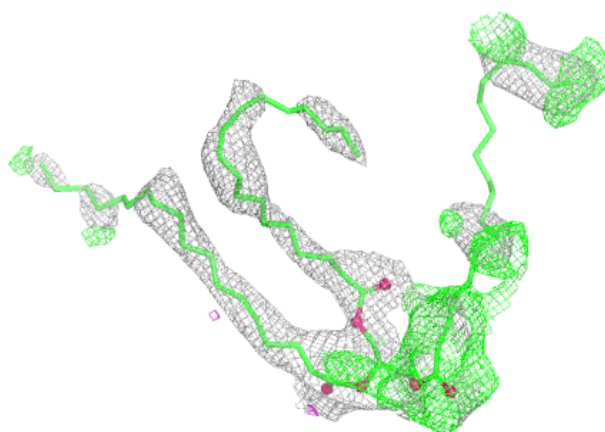


Electron density around CDL G 102:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

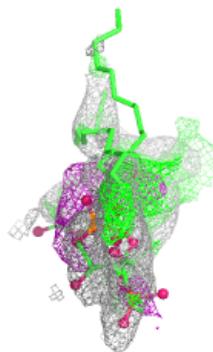
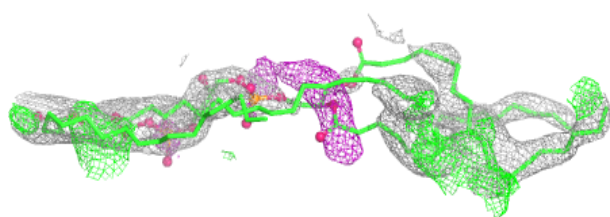
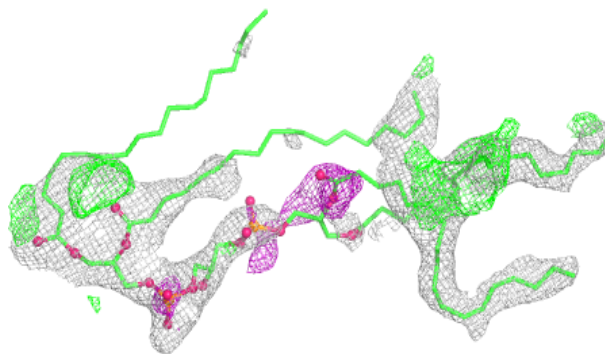
**Electron density around TGL D 201:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



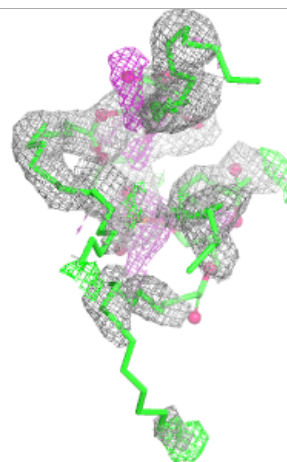
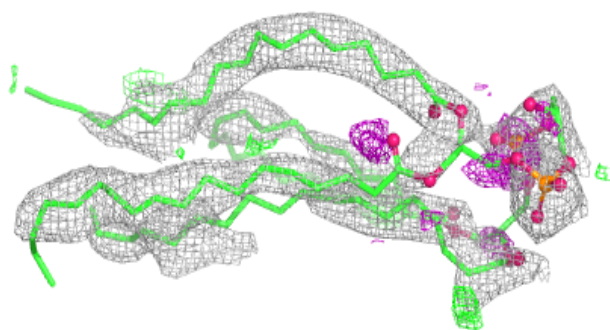
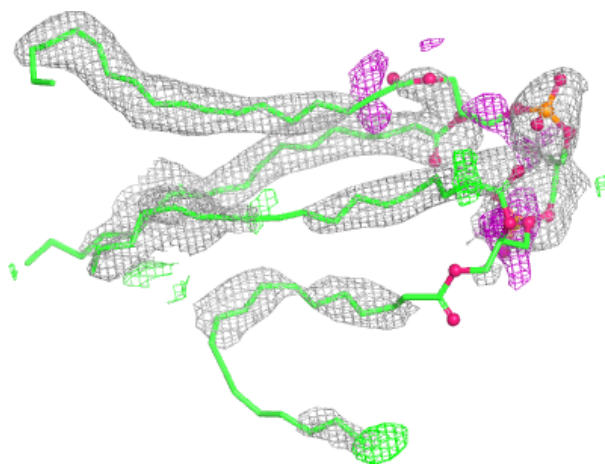
Electron density around CDL T 104:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



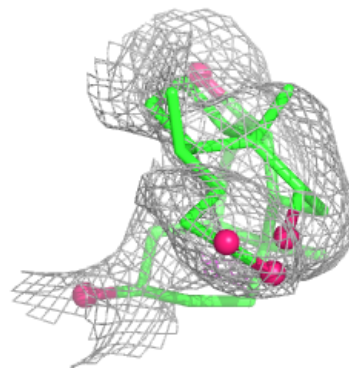
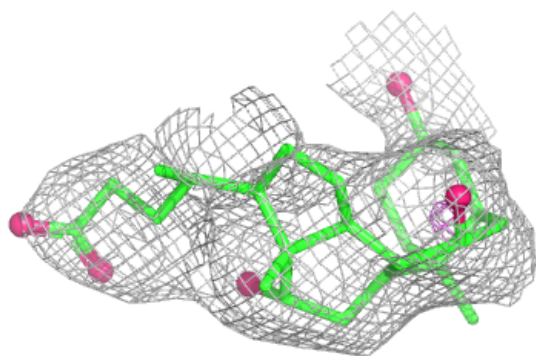
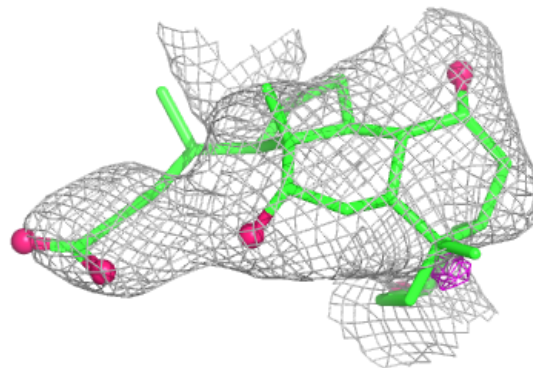
Electron density around CDL P 304:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



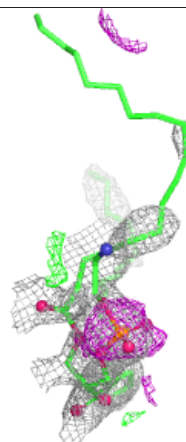
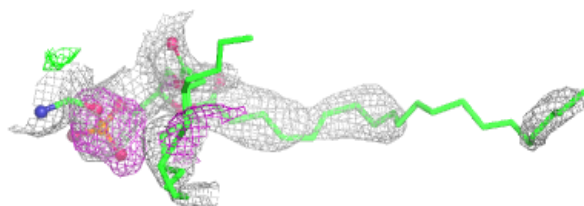
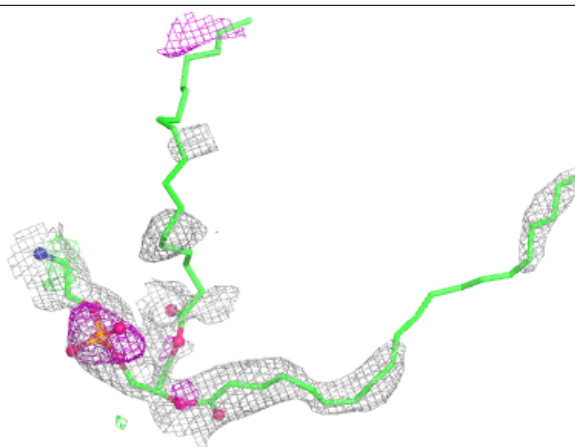
Electron density around CHD W 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



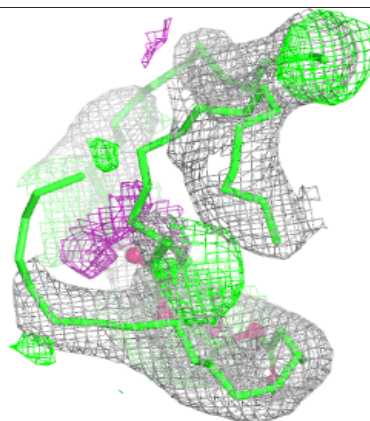
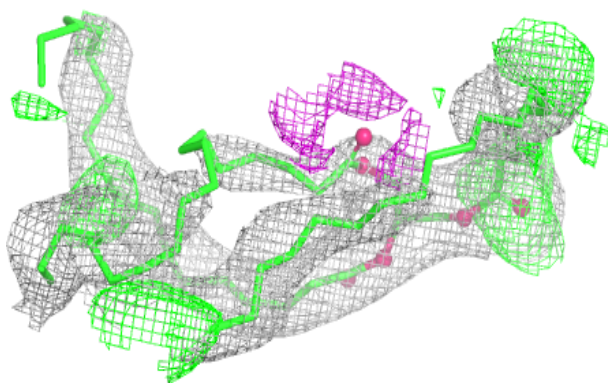
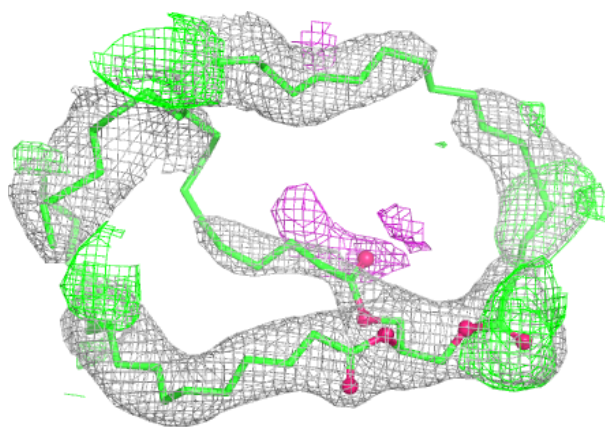
Electron density around PEK G 103:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



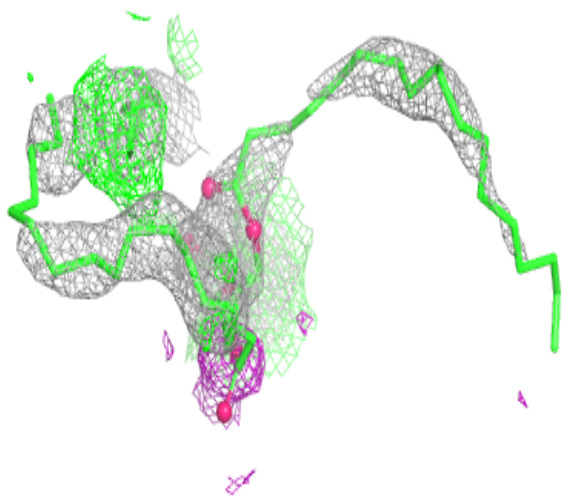
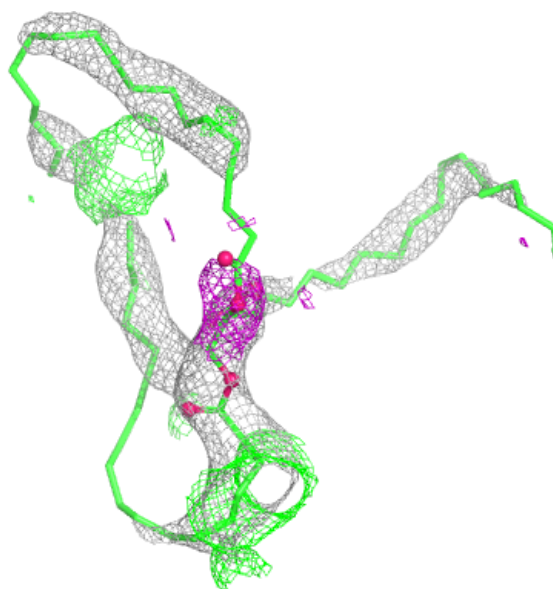
Electron density around TGL A 607:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



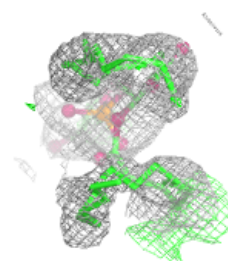
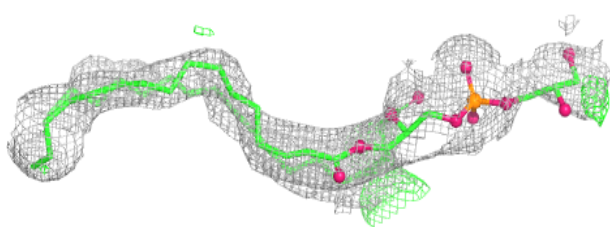
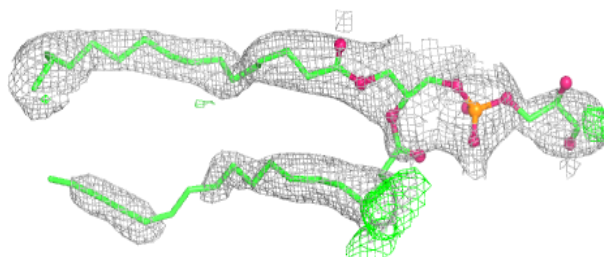
Electron density around TGL Y 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

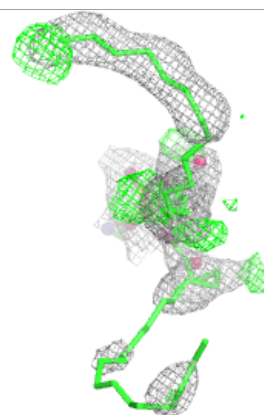
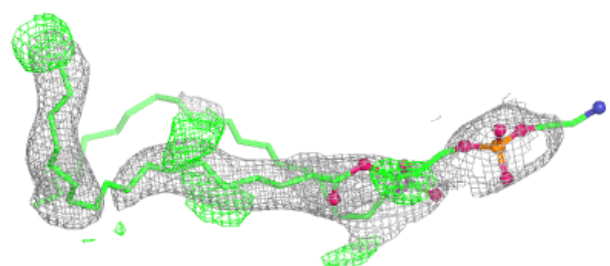
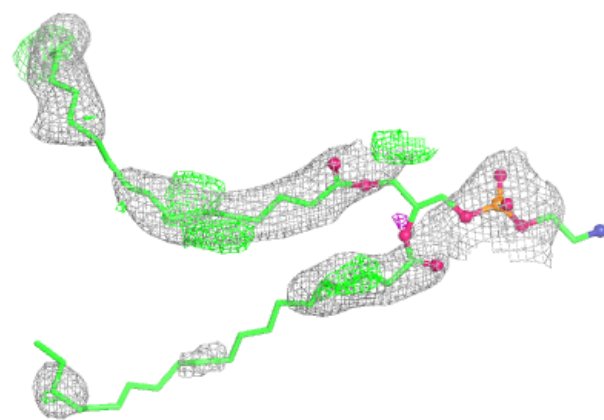


Electron density around PGV P 301:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

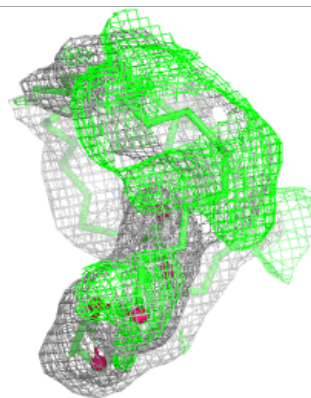
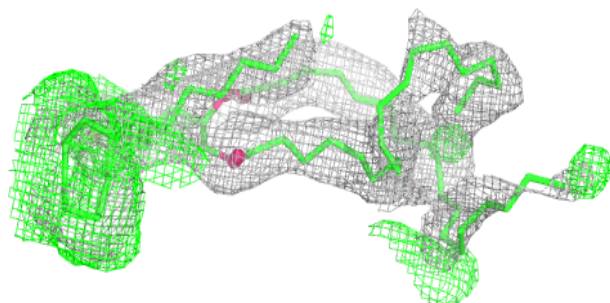
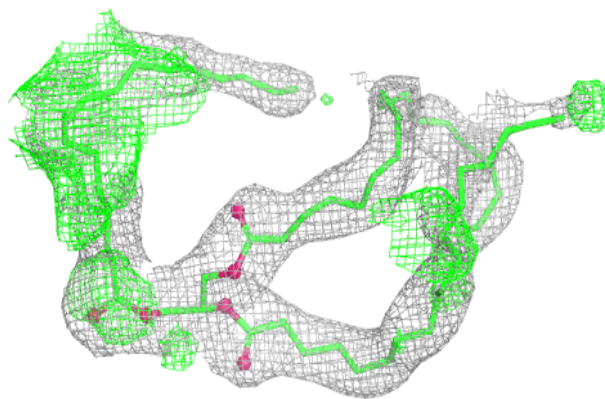
**Electron density around PEK C 306:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

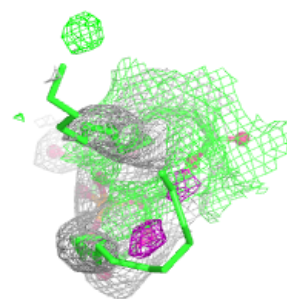
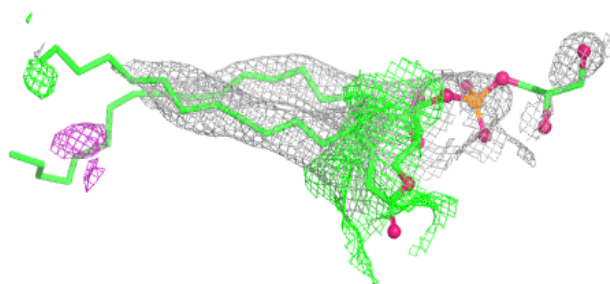
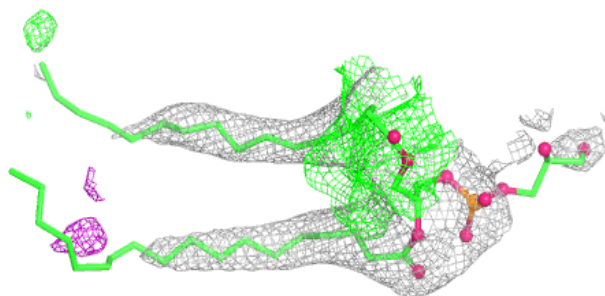


Electron density around TGL O 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

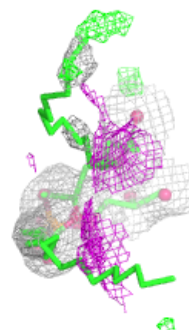
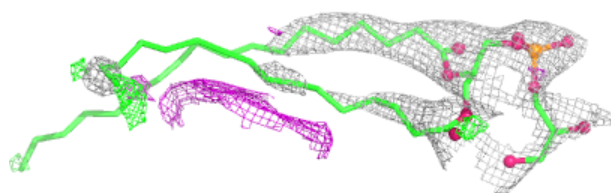
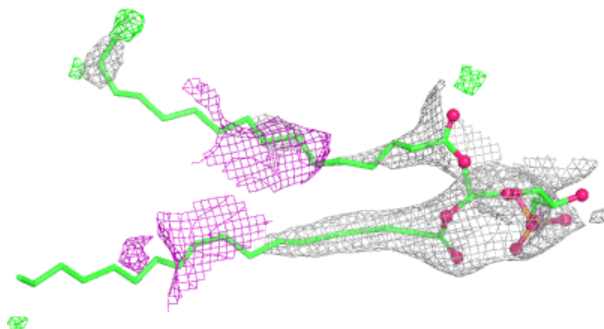
**Electron density around PGV A 609:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

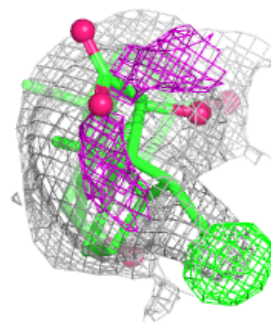
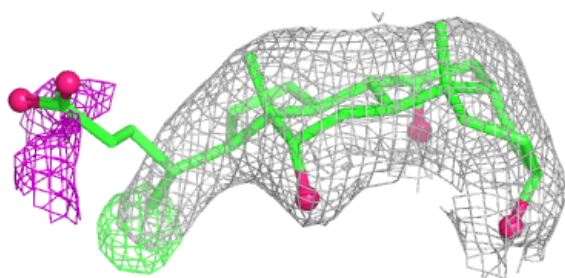
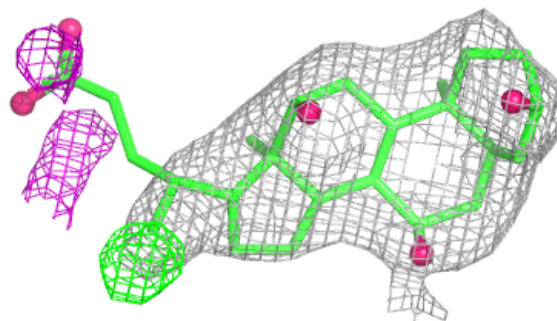


Electron density around PGV N 607:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

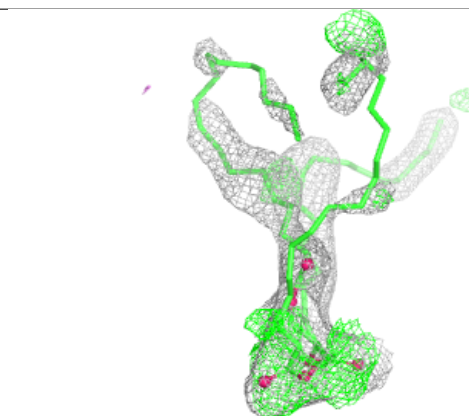
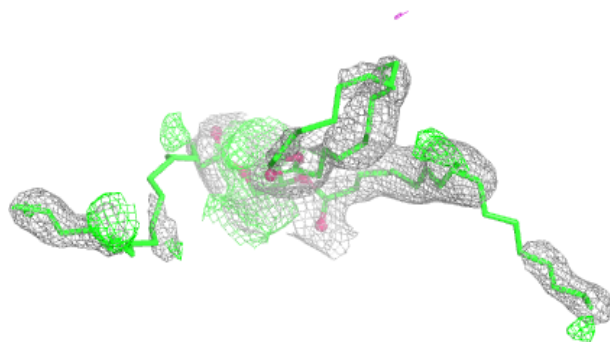
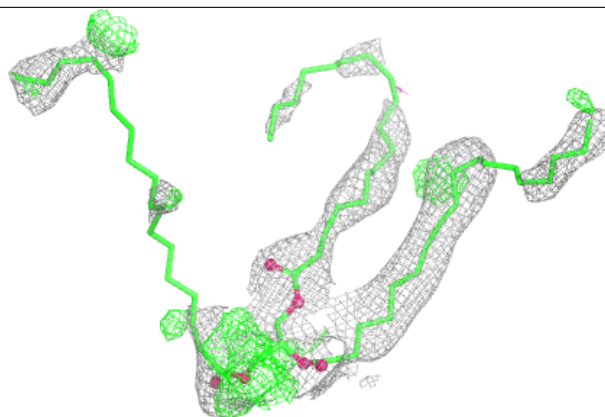
**Electron density around CHD J 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

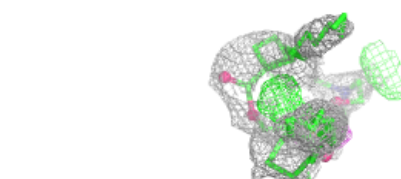
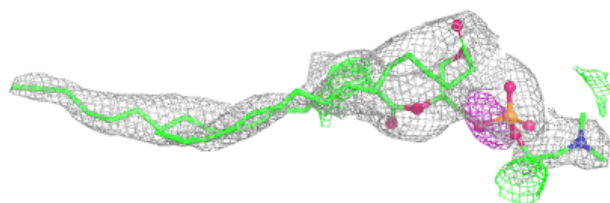
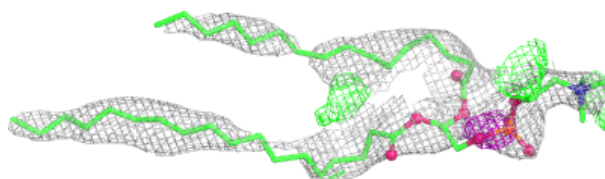


Electron density around TGL Q 201:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

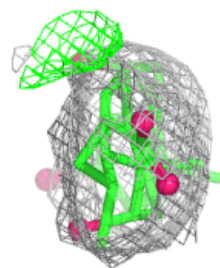
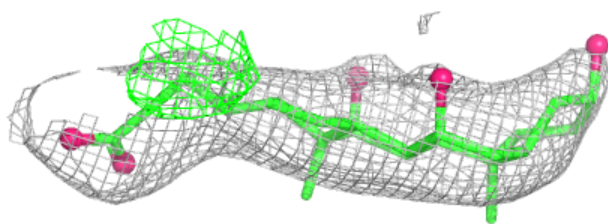
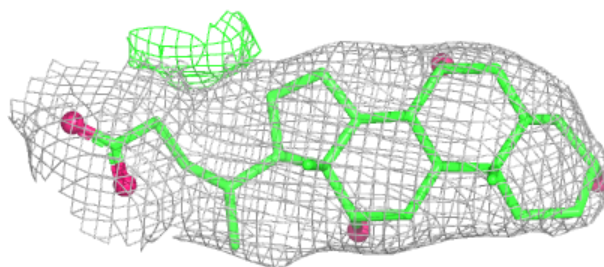
**Electron density around PSC B 303:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

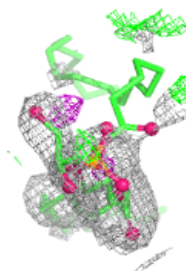
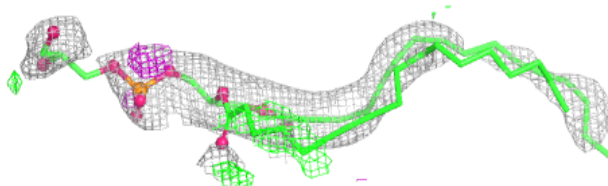
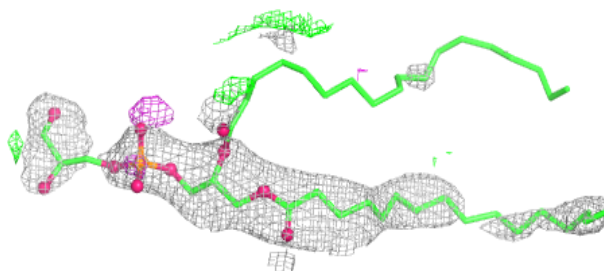


Electron density around CHD P 305:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

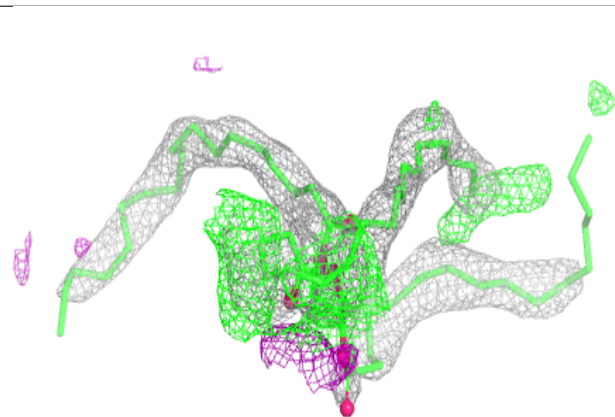
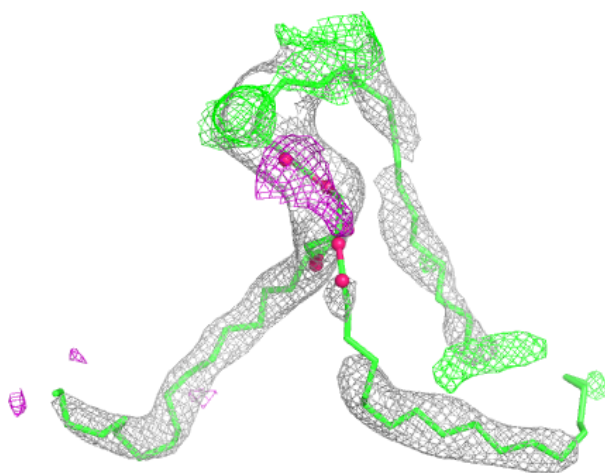
**Electron density around PGV C 307:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



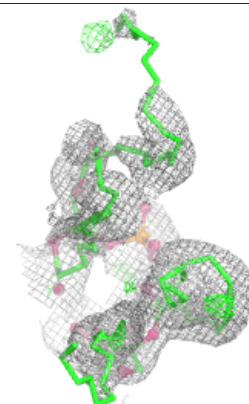
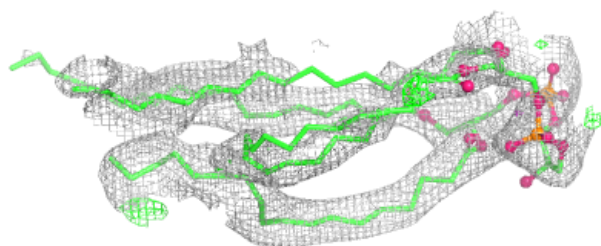
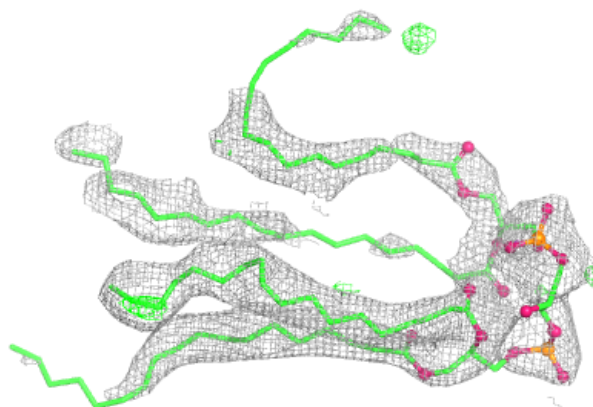
Electron density around TGL L 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

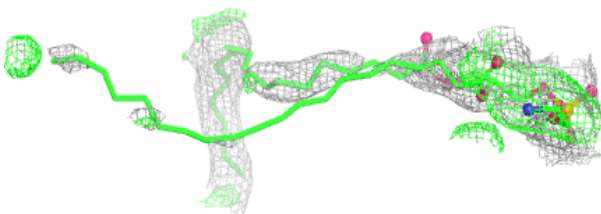
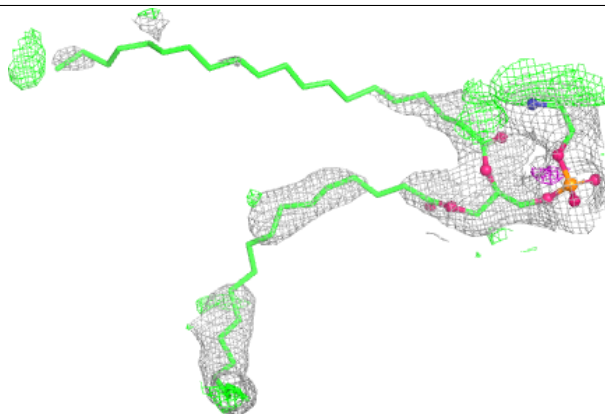


Electron density around CDL C 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

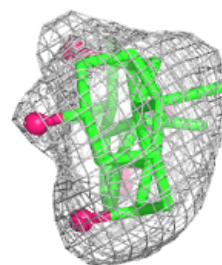
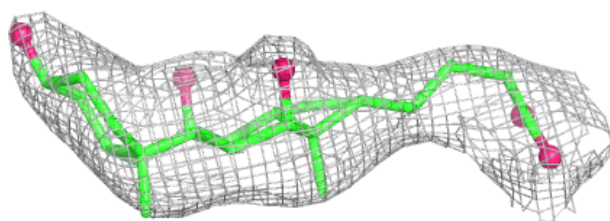
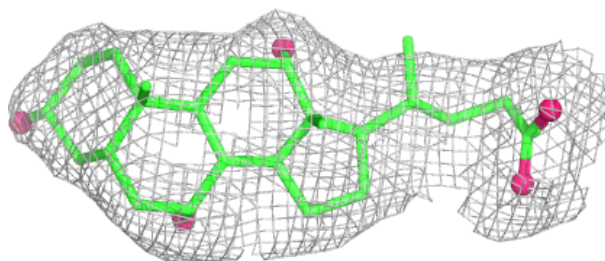
**Electron density around PEK T 103:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

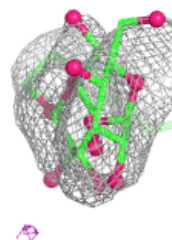
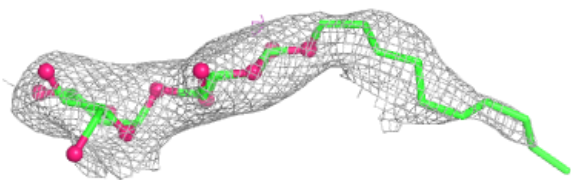
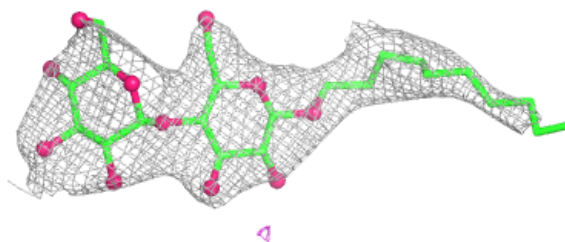


Electron density around CHD C 304:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

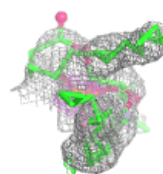
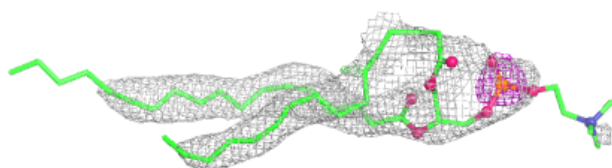
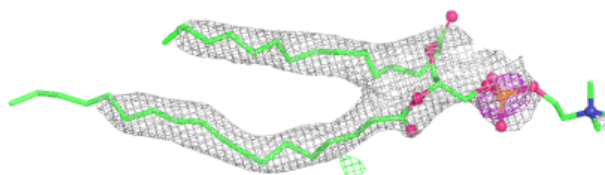
**Electron density around DMU Z 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

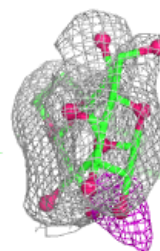
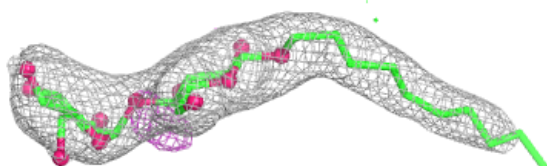
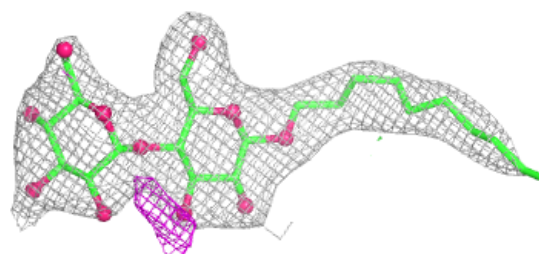


Electron density around PSC O 303:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

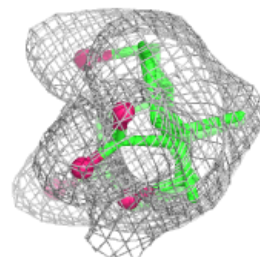
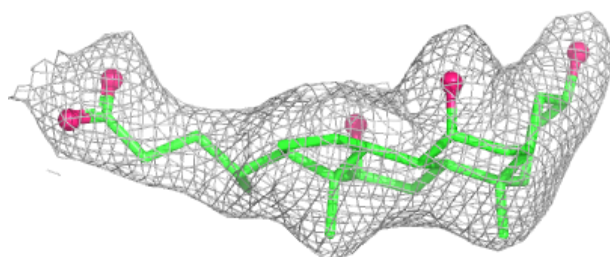
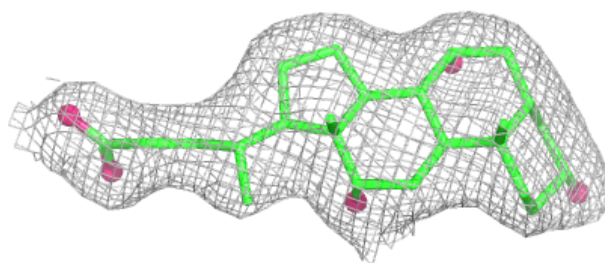
**Electron density around DMU M 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

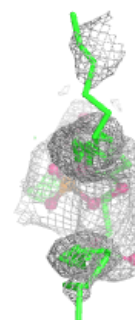
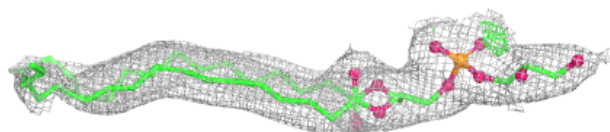
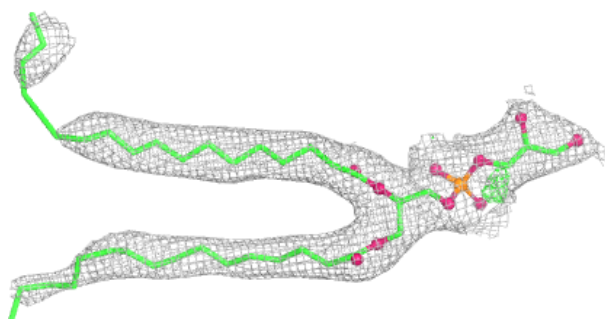


Electron density around CHD P 306:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

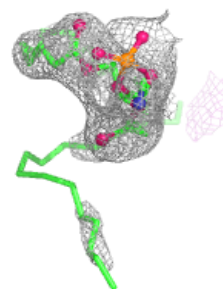
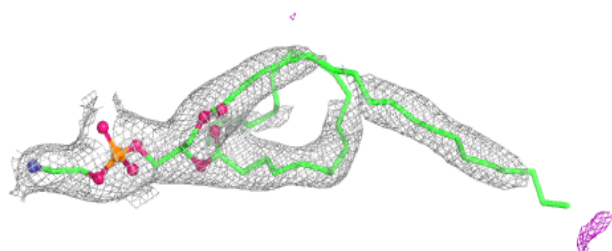
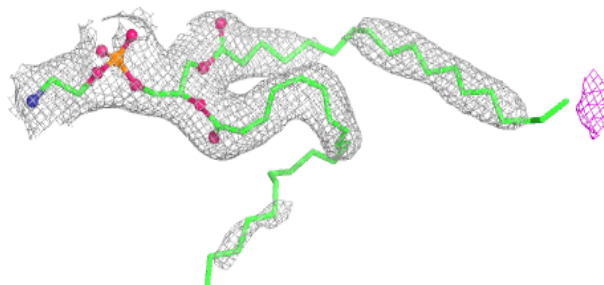
**Electron density around PGV P 303:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

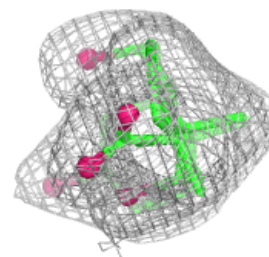
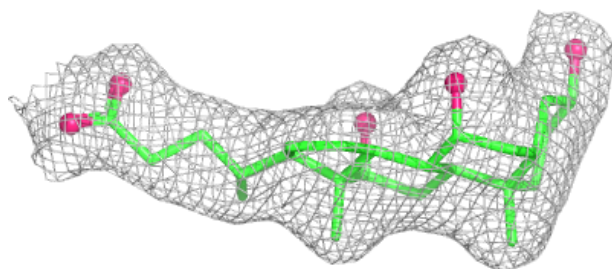
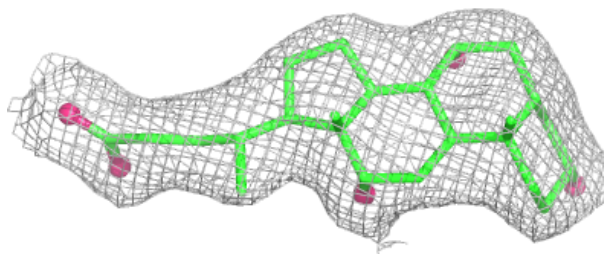


Electron density around PEK T 101:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

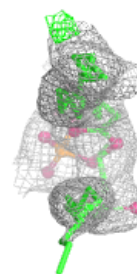
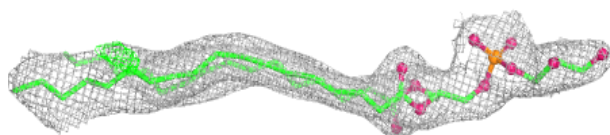
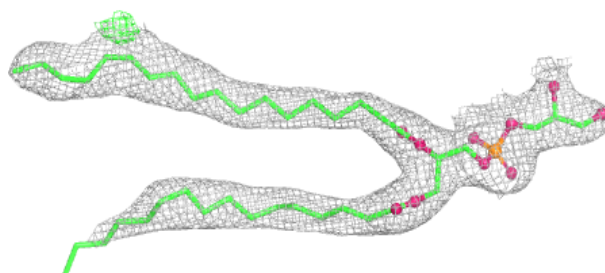
**Electron density around CHD C 305:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

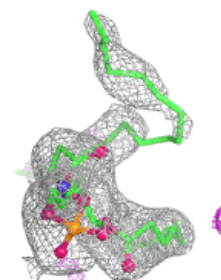
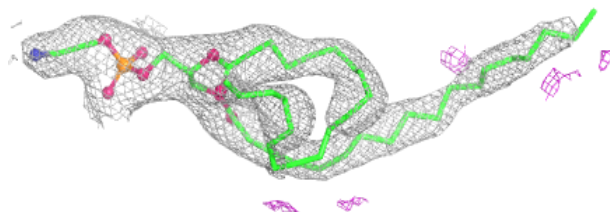
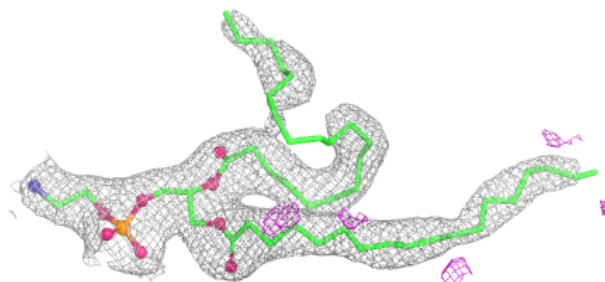


Electron density around PGV C 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

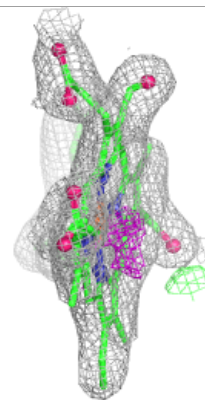
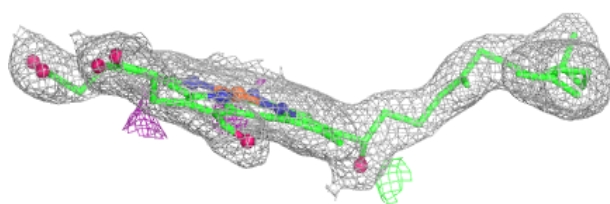
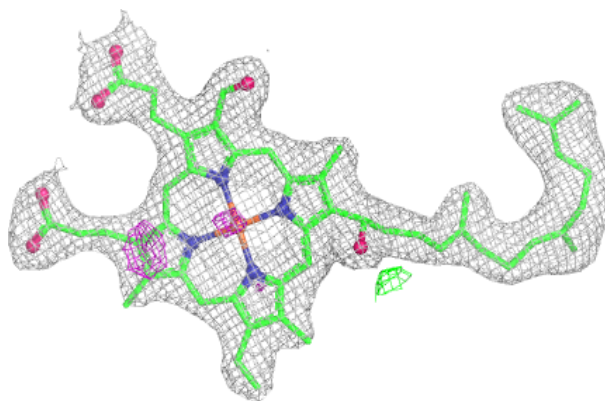
**Electron density around PEK G 101:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

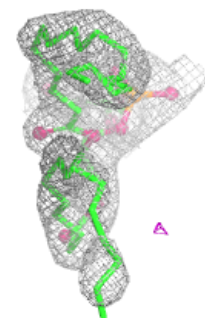
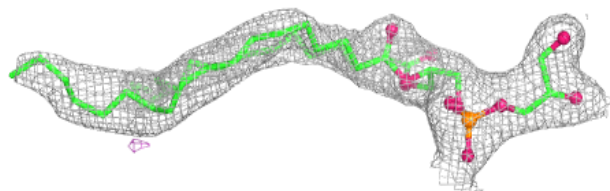
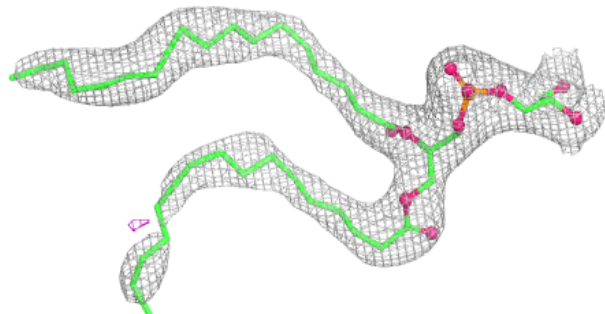


Electron density around HEA A 602:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

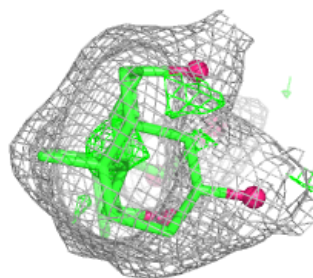
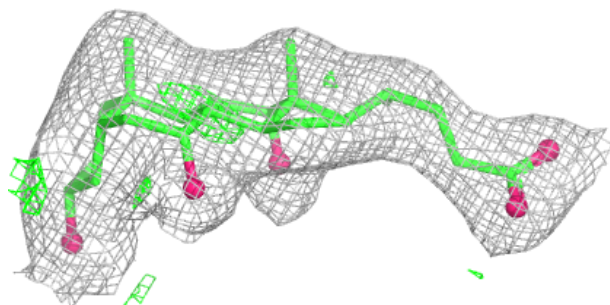
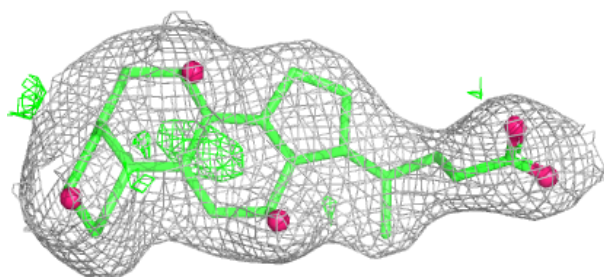
**Electron density around PGV N 608:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

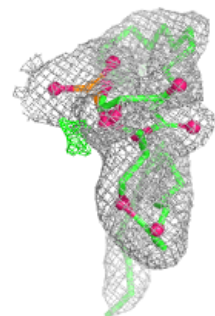
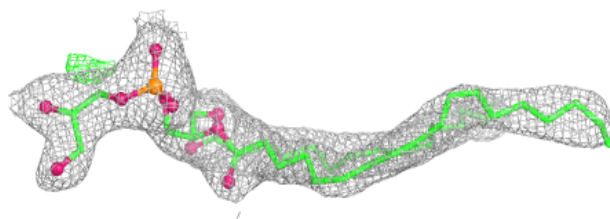
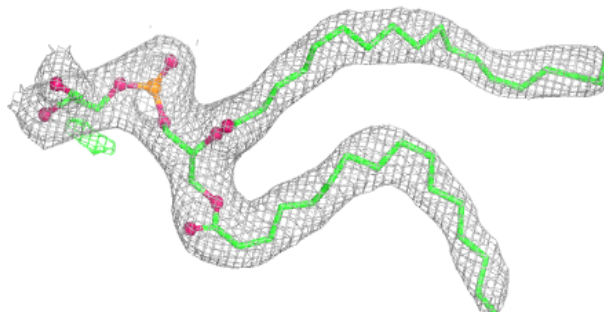


Electron density around CHD G 104:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

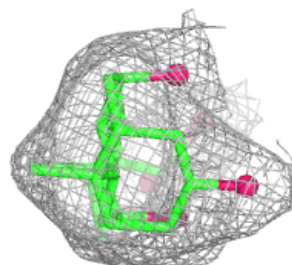
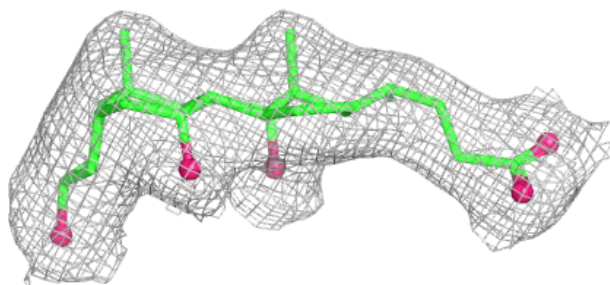
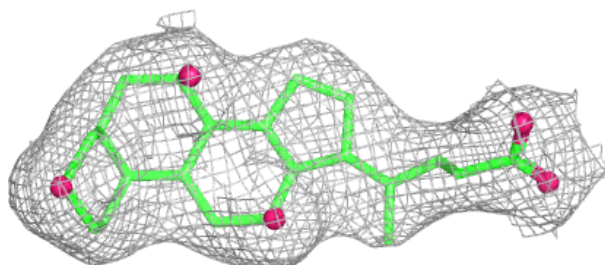
**Electron density around PGV A 608:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

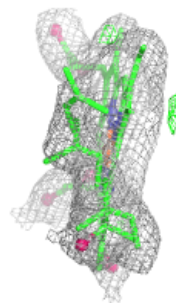
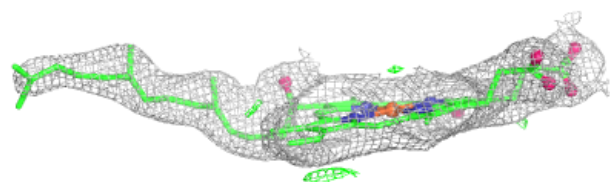
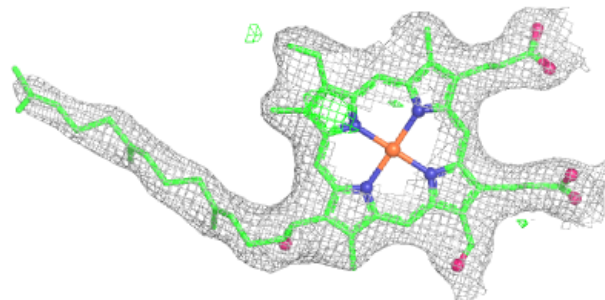


Electron density around CHD B 302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

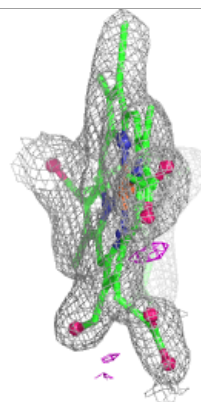
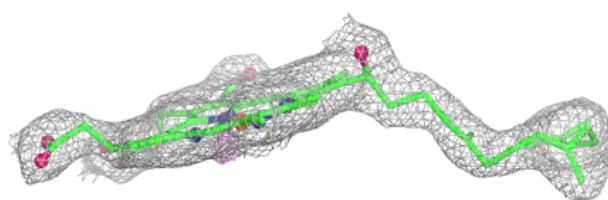
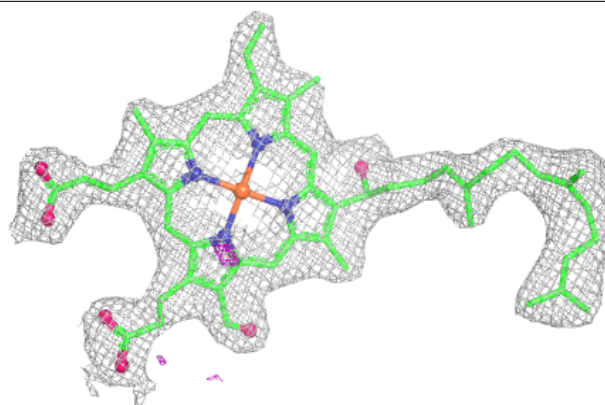
**Electron density around HEA N 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

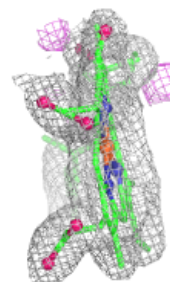
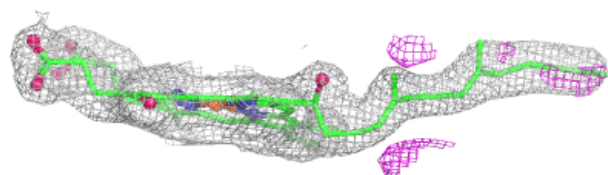
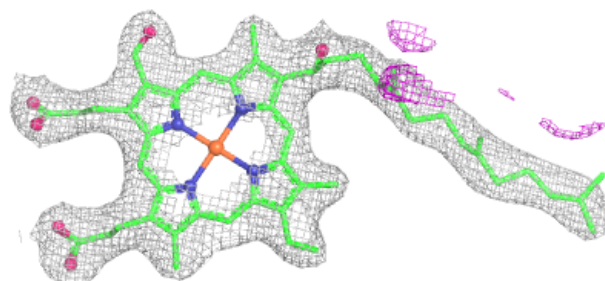


Electron density around HEA N 602:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around HEA A 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.